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SOME RESULTS IN THE THEORY OF SUBSET SELECTION PROCEDURES, (U)

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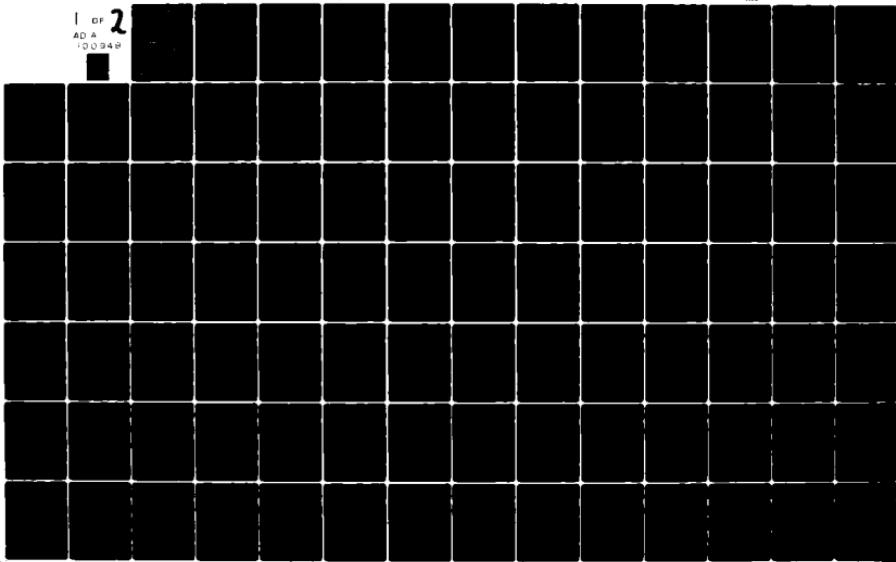
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Some Results in the Theory  
of Subset Selection Procedures

by

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INTRODUCTION

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Selection and ranking (ordering) problems in statistical inference arise mainly because the classical tests of homogeneity are often inadequate in certain situations where the experimenter is interested in comparing  $k ( \geq 2 )$  populations, treatments or processes with the goal of selecting one or more worthwhile (good) populations. Mosteller (1949), Paulson (1949), Bahadur (1950) and Bahadur and Robbins (1950) were among the earliest research workers to recognize this inadequacy and to formulate the problem as a multiple decision problem aimed at the selection and ranking of the  $k$  populations.

In the thirty years since these early papers, selection and ranking problems have become an active area of statistical research. There have been two approaches to these problems, the 'indifference zone' approach and the 'subset selection' approach. In the first approach, due to Bechhofer (1954), the experimenter wishes to select one population (or a fixed number  $t \geq 1$  of population) which is guaranteed to be the one of interest to him with a fixed probability  $P^*$  whenever the unknown parameters lie outside some subspace of the parameter space, the so-called indifference zone. Important contributions using this approach have been made by Bechhofer and Sobel (1954), Bechhofer, Dunnett and Sobel (1954), Sobel and Huyett (1957), Sobel (1962), Bechhofer, Trotter and Sobel (1968), Mahanunulu (1967), Desu and Sobel (1962, 1971) and

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Tamhane and Bechhofer (1977, 1979) among others. A quite complete bibliography may be found in Gupta and Panchapakesan (1979) (see also Gibbons, Olkin and Sobel (1977)).

The second approach pioneered by Gupta (1956, 1963, 1965) assumes no a priori information about the parameter space. A single population is not necessarily chosen; rather a subset of the given  $k$  populations is selected depending on the outcome of the experiment. It is guaranteed to contain the population(s) of interest with probability which is at least equal to  $P^*$  (the basic probability requirement) regardless of the true unknown configurations of the parameters. Some recent contributors in the category of subset selection include: Beely (1965), Gnanadesikan (1966), Gnanadesikan and Gupta (1970), Gupta (1967), Gupta and Studden (1970), Nagel (1970), Gupta and Nagel (1971), Gupta and Panchapakesan (1972), Rizvi and Sobel (1967), McDonald (1969), Gupta and McDonald (1970), Santner (1975), W. T. Huang (1972), D. Y. Huang (1975), Gupta and Huang (1975a, 1975b) and Gupta and Huang (1976).

Subset selection procedures can also be thought of as screening procedures which enable the experimenter to select a subset of populations (under study) which contains the populations of interest so that the populations in the selected subset can be further studies.

Sequential and multistage aspects of the ranking and selection problems, have been explored, based on the indifference zone approach by Bechhofer, Dunnett and Sobel (1954), Bechhofer (1958), Paulson (1962, 1963, 1964, 1967) and Bechhofer, Fiefer and Sobel (1968). Barson and Gupta (1972), Huang (1972), Gupta and Huang (1975), Gupta and Miescke (1979) and Carroll (1974) have investigated subset selection procedures, based on sequential sampling.

Contributions to optimum properties of subset selection procedures have been made by Lehmann (1961), Studden (1967), Deely and Gupta (1968), Berger (1977, 1979), Gupta and Hsu (1978), Gupta and Miescke (1978), Berger and Gupta (1980).

In the decision-theoretic approach to the subset selection problem, Goel and Rubin (1977), Chernoff and Yahav (1977), Bickel and Yahav (1977), Gupta and Hsu (1978), Miescke (1979), Gupta and Kim (1980), Gupta and Hsiao (1980) have given different formulations under different loss functions and carried out investigations which indicate that the gamma type maximum (minimum) means procedures are quite 'optimal' and 'robust'.

The main purpose of this thesis is to study some problems using the subset selection approach and provide procedures and results for some unsolved problems.

Chapter I considers the problem of selecting a subset containing all populations better than a control under an ordering prior. Here, by an ordering prior we mean that there exists a known simple or partial order relationship among the unknown parameters of the treatments (excluding the control). Three new selection procedures are proposed and studied. These procedures do meet the usual requirement that the probability of a correct selection is greater than or equal to a pre-determined number  $P^*$ . Two of the three procedures use the isotonic regression over the sample means of the  $k$  treatments with respect to (wrt) the given ordering prior. Tables which are necessary to carry out the selection procedures with isotonic approach for the selection of unknown means of normal populations and gamma populations are given. Monte Carlo comparisons on the performance of several procedures for the normal or gamma means problem were carried out in

several selected cases; these are given in Table V and Table VI at the end of Chapter I. In each case ten thousand simulations were performed. The results of this study seem to indicate that the procedures based on isotonic estimators always have superior performance, especially, when there are more than one bad populations (in comparison with the control).

Chapter II deals with a new 'Bayes- $P^*$ ' approach about the problem of selecting a subset which contains the 'best' of  $k$  populations. Here, by best we mean the (unknown) population with the largest unknown mean. The (non-randomized) Bayes- $P^*$  rule refers to a rule with minimum risk in the class of (non-randomized) rules which satisfy the condition that the posterior probability of selecting the best is at least equal to  $P^*$ . Given the priors of the unknown parameters, two 'Bayes- $P^*$ ' subset selection procedures  $\psi^B$  and  $\psi_{NR}^B$  (randomized and non-randomized, respectively) under certain loss functions are obtained and compared with the classical maximum-type means procedure  $\psi^M$ . The comparisons of the performance of  $\psi^B$  with  $\psi_{NR}^B$  and  $\psi^M$ , based on Monte Carlo studies, indicate that the procedure  $\psi^B$  always has higher 'efficiency' and smaller expected selected size of the selected subset. Also  $\psi^B$  appears to be robust when the true distributions are not normal but are some other symmetric distributions such as, the logistic, the double exponential, Laplace, and the gross error model (the contaminated distribution).

CHAPTER I  
SELECTION PROCEDURES FOR POPULATIONS  
BETTER THAN A CONTROL UNDER ORDERING PRIOR

1.1. Introduction

In this chapter, three new selection procedures are given for the problem of selecting a subset which contains all populations better than a standard or control under simple or partial ordering prior. Here by simple or partial ordering prior we mean that there exist known simple or partial order relationships (defined more specifically later in Section 1.2) among unknown parameters. The procedures described do meet the usual requirement that the probabilities of a correct selection are greater than or equal to a predetermined number  $P^*$ , the so-called  $P^*$  condition.

Many authors have considered the problem of comparing populations with a control under different types of formulations (see Gupta and Panchapakesan (1979)). Dunnett (1955) considered the problem of separating those treatments which are better than the control from those that are worse. Gupta and Sobel (1958), Gupta (1965), Naik (1975), Broström (1977) studied the problem of selecting a subset containing all populations better than the control. Lehmann (1961) discussed similar problems with emphasis on the derivation of a restricted minimax procedure. Kim (1979), Hsiao (1979) studied the problem of

selecting populations close to a control. In all these papers it is assumed that all populations are independent and that there is no information about the order of unknown parameters. However, in many situations, we may know something about the unknown parameters. What we know is always not the prior distributions but some partial or incomplete prior information, such as the simple or partial order relationship among the unknown parameters. This type of information about the ordering prior may come from the past experiences; or it may arise in the experiments where, for example, higher dose level of some drugs always has larger effect (side-effect) on the patients.

In Section 1.2 definitions and notations used in this chapter are introduced. In Section 1.3 we consider the problem for location parameters. We propose three types of selection procedures for the cases when the control parameter is known or not known (the scale parameter may or may not be assumed known). Some equivalent forms of the procedures are given, and their properties are discussed. In Section 1.4 the problem for scale parameters of the gamma distributions is considered and three analogous selection procedures are proposed. In both Section 1.3 and 1.4 simple ordering priors are assumed and some theorems in the theory of random walks are used. In Section 1.5 a selection procedure is given for the problem of selecting all populations better than the control under partial ordering prior. Section 1.6 deals with the use of Monte Carlo techniques to make comparisons among the selection procedures proposed in Section 1.3 and those in Section 1.4, respectively.

## 1.2. Notations and Definitions

Suppose we have  $k + 1$  populations  $\pi_0, \pi_1, \dots, \pi_k$ . The population treatment  $\pi_0$  is called the control or standard population. Assume that the random variables  $X_{ij}$  associated with  $F(\cdot; \pi_i)$  and  $\bar{X}_{ij}, \dots, \bar{X}_{in_i}$ ,  $i = 1, \dots, k$ , is an independent sample from  $\pi_i$ . Assume that we have an ordering prior of  $\pi_1, \dots, \pi_k$ . First we assume that the ordering prior is the simple order, so that without loss of generality, we may assume that,  $\pi_1 \leq \dots \leq \pi_k$ . In Section 1.5 we will consider the partial ordering prior case. Note that the values of  $\pi_i$ 's are unknown.

Suppose our goal is to find a non-trivial (small) subset which contains all populations with parameter larger (smaller) than the control  $\pi_0$  (known or unknown) with probability not less than a given value  $\pi^*$ .

The action space  $\mathcal{C}$  is the class of all subsets of set  $\{1, 2, \dots, k\}$ . An action  $A$  is the selection of some subset of the  $k$  populations. This means that  $\pi_i$  is included in the selected subset.

Let  $\pi = (\pi_0, \pi_1, \dots, \pi_k)$ . Then the parameter space is denoted by  $\mathcal{S}$ , where  $\mathcal{S} = \{\pi \in \mathbb{R}^{k+1} \mid \pi_1 \leq \pi_2 \leq \dots \leq \pi_k; \pi \leq \pi_0 \leq \pi\}$  is a subset of  $k + 1$  dimensional Euclidean space  $\mathbb{R}^{k+1}$ .

The sample space is denoted by  $\mathcal{X}$  where

$$\mathcal{X} = \{x \in \mathbb{R}^{n_1 + \dots + n_k} \mid x = (x_{11}, \dots, x_{1n_1}, x_{21}, \dots, x_{kn_k}, \dots, x_{in_i})\}.$$

**Definition 1.2.1.** A (non-randomized) selection procedure (rule)  $\pi^*$  is a mapping from  $\mathcal{S}$  to  $\mathcal{C}$ .

A population  $\pi_i$  ( $i = 1, \dots, k$ ) is called a good population if  $\pi_i \geq \pi_0$ , and we say a selection procedure  $\delta$  make a correct selection (CS) if the selected subset contains all good populations. A selection procedure  $\delta$  satisfies the  $P^*$ -condition if

$$P_\delta(\text{CS}|\delta) \geq P^* \text{ for all } \delta \in \Delta$$

that is

$$\inf_{\delta \in \Delta} P_\delta(\text{CS}|\delta) \geq P^*. \quad (1.2.1)$$

Let  $\Delta = \{\delta \mid \inf_{\delta \in \Delta} P_\delta(\text{CS}|\delta) \geq P^*\}$  be a collection of all selection procedures satisfying the  $P^*$ -condition.

In the sequel we will use the isotonic estimators (see Barlow, Bartholomew, Bremner and Brunk (1972)). Hence we give the following definitions and theorems.

**Definition 1.2.2.** Let the set  $\mathcal{S}$  be a finite set. A binary relation " $\leq$ " on  $\mathcal{S}$  is called a simple order if it is

- (1) reflexive:  $x \leq x$  for  $x \in \mathcal{S}$
- (2) transitive:  $x, y, z \in \mathcal{S}$  and  $x \leq y, y \leq z$  imply  $x \leq z$
- (3) antisymmetric:  $x, y \in \mathcal{S}$  and  $x \leq y, y \leq x$  imply  $x = y$
- (4) every two elements are comparable:  $x, y \in \mathcal{S}$  imply either  $x \leq y$  or  $y \leq x$ .

A partial order on  $\mathcal{S}$  is a binary relation " $\leq$ " on  $\mathcal{S}$ , such that it is (1) reflexive, (2) transitive, and (3) antisymmetric. Thus every simple order is a partial order. We use poset  $(\mathcal{S}, \leq)$  to denote the set  $\mathcal{S}$  that has a partial order binary relation " $\leq$ " on it.

**Definition 1.2.3.** A real-valued function  $f$  is called isotonic on poset  $(\mathcal{S}, \leq)$  if and only if (1)  $f$  is defined on  $\mathcal{S}$ , (2) if  $x, y \in \mathcal{S}$  and  $x \leq y$  then  $f(x) \leq f(y)$ .

**Definition 1.2.4.** Let  $q$  be a real-valued function on  $\mathcal{S}$  and let  $W$  be a given positive function on  $\mathcal{S}$ . A function  $q^*$  on  $\mathcal{S}$  is called an isotonic regression of  $q$  with weights  $W$  if and only if:

(1)  $q^*$  is an isotonic function on poset  $(\mathcal{S}, \leq)$

$$(2) \quad \sum_{x \in \mathcal{S}} [q(x) - q^*(x)]^2 W(x) = \min_{f \in \mathcal{Q}} \sum_{x \in \mathcal{S}} [q(x) - f(x)]^2 W(x),$$

where  $\mathcal{Q}$  is the class of all isotonic functions on poset  $(\mathcal{S}, \leq)$ .

From Barlow, et. al. (1972), (see their Theorem 1, Cor. 1, and the corollary there), we have the following theorem.

**Theorem 1.2.1.** There exists one and only one isotonic regression  $q^*$  of  $q$  with weight  $W$  on poset  $(\mathcal{S}, \leq)$ .

**Definition 1.2.5.** A set  $S$  is convex if  $s_1, s_2 \in S$  and if  $0 \leq t \leq 1$  then  $s_1 + (1-t)s_2 \in S$ .

**Definition 1.2.6.** A set  $S$  is a cone if  $s \in S$  then for any non-negative real number  $c$ ,  $cs \in S$ .

**Definition 1.2.7.** A poset  $(\mathcal{S}, \leq)$  is a lattice if for all  $A$  and  $B$  in  $\mathcal{S}$  there exist for any finite non-empty subset  $U$  of  $\mathcal{S}$ ,

If  $f$  and  $g$  are two isotonic functions on poset  $(\mathcal{S}, \leq)$ , we define  $f \wedge g$  and  $f \vee g$  as

$$(f \wedge g)(t) = f(t) \wedge g(t) = \min(f(t), g(t))$$

and

$$(f \vee g)(t) = r(t) \vee g(t) = \max(r(t), g(t)).$$

Then we state the following:

**Theorem 1.2.2.** The class  $\mathcal{E}$  of all isotonic functions on poset  $(\mathcal{S}, \leq)$  is a convex cone and a lattice.

There are some algorithms, such as the "pool-adjacent-violation" algorithm (see page 13 of Barlow, et. al. (1972)) or Ayer, Frank, Tukey, Reid and Silverman (1955) or the "up-and-down blocks" algorithm, see, e.g. (1964), which show how to calculate the isotonic regression under simple order.

The following max-min formulas were given by Ayer et. al. (1955).

**Theorem 1.2.3. (max-min formulas)**

Assume that we have poset  $(\mathcal{S}, \leq)$  where  $\mathcal{S} = \{1, \dots, k\} \cup \{\infty\}$  and that function  $g: \mathcal{S} \rightarrow \mathbb{R}$ , then the isotonic regression  $g^*$  of  $g$  with weight  $W$  has the following formulas:

$$\begin{aligned} g^*(i) &= \max_{s \leq i \leq t} \text{Av}(s, t) \\ &= \max_{s \leq i \leq t} \text{Av}(s, t) \\ &= \min_{t \leq i \leq s} \max_{s \leq t} \text{Av}(s, t) \\ &= \min_{t \leq i \leq s} \max_{s \leq t} \text{Av}(s, t) \end{aligned}$$

where

$$\text{Av}(s, t) = \frac{\sum_{r=s}^t w(r)}{\sum_{r=s}^t w(r)}.$$

Corollary 1.2.1.  $(q + c)^* = q^* + c$ ,

$$(aq)^* = aq^*, \text{ if } a > 0.$$

Corollary 1.2.2.  $[(\eta^*)q + \zeta(\eta^*)]^* = (\eta^*)q^* + \zeta(\eta^*)$ , where  $\eta^*$  is a non-negative function and  $\zeta$  is an arbitrary function.

### 1.3. Proposed Selection Procedures for the Location Parameter Problem

To discuss some more general results, we assume that population  $\eta$  has an absolutely continuous location-scale distribution function

$F(x; \eta_1, \eta_2) = F\left(\frac{x-\eta_1}{\eta_2}\right)$ , where  $0 < F(x) < 1$  for all  $x$ ,  $\eta_1 < \eta_2$  and the values of  $\eta_1, \dots, \eta_k$  are unknown, but their ordering, say,  $\eta_1 < \dots < \eta_k$  is known. Note that in this case we replace  $\eta$  in the parameter space by  $\eta$ , all other quantities remaining the same.

Let us define the subspace  $\eta = (\eta_1, \eta_2, \dots, \eta_k) \in \eta_0 = \{(\eta_1, \eta_2, \dots, \eta_k) \mid \eta_1 < \eta_2 < \dots < \eta_k\}$  for  $i = 1, \dots, k-1$  and let subspace  $\eta_k = (\eta_k) \in \eta_0 - \eta_1$  and subspace

$\eta_0 = \{(\eta_1, \eta_2, \dots, \eta_k) \in \eta_0\}$ , then we have  $\eta = \sum_{i=0}^k \eta_i$ . Note that the constant  $\eta_0$  could be known or unknown. If  $\eta_0$  is unknown, we assume that the distribution of population  $\eta_0$  is  $F(x; \eta_0, \eta_1)$  and we take independent observations  $x_{01}, \dots, x_{0n_0}$  from  $\eta_0$  and the sample space  $\Omega$  turns to

$\{x \in \mathbb{R}^{n_0 + \dots + n_k} \mid x = (x_{01}, \dots, x_{0n_0}, x_{11}, \dots, x_{1n_1}, \dots, x_{kn_k})\}$ . Using the partition  $\eta_0, \dots, \eta_k$  of parameter space  $\eta$ , we have

$$\inf_{\eta \in \Omega} P_{\eta}(\text{CS}^{\star}) = \inf_{\eta_0 \in \Omega_0} \inf_{\eta_1 \in \Omega_1} \dots \inf_{\eta_k \in \Omega_k} P_{\eta}(\text{CS}^{\star}),$$

for any selection procedure  $\text{CS}^{\star}$ . Hence the  $P^{\star}$ -condition is equivalent to

$$\inf_{\mu \in \mathcal{C}_i} P_{\mu}(CS|s) \geq P^*, \text{ for } i = 1, \dots, k.$$

Note that  $\inf_{\mu \in \mathcal{C}_0} P_{\mu}(CS|s) = 1$  for any selection procedure since there

exist no good population in this case.

Suppose  $\bar{x}_i = x_i$  is the outcome of the sample mean of population  $\mu_i$ ,  $i = 1, \dots, k$ . Let  $\mathcal{S}$  denote the set  $\{u_1, u_2, \dots, u_k\}$  where  $u_1 \leq \dots \leq u_k$ , and let  $w(u_i) = n_i^{-1/2} - w_i$ ,  $g(u_i) = x_i$ ,  $i = 1, \dots, k$ . Then by the max-min formulas, the isotonic regression of  $g$  is  $g^*$ , where

$$g^*(u_i) = \max_{1 \leq s \leq i} \min_{s \leq t \leq k} \frac{\sum_{j=s}^t x_j w_j}{\sum_{j=s}^t w_j}, \quad i = 1, \dots, k.$$

The isotonic estimator of  $\mu_i$  is denoted by  $\hat{x}_{i:k}$ ,  $i = 1, \dots, k$  where

$$\begin{aligned} \hat{x}_{i:k} &= \max_{1 \leq s \leq i} \min_{s \leq t \leq k} \frac{\sum_{j=s}^t x_j w_j}{\sum_{j=s}^t w_j} \\ &= \max_{1 \leq j \leq i} \hat{x}_{j:k} \end{aligned} \quad (1.3.1)$$

where

$$\hat{x}_{j:k} = \min \left( x_j, \frac{x_j w_j + x_{j+1} w_{j+1}}{w_j + w_{j+1}}, \dots, \frac{x_j w_j + \dots + x_k w_k}{w_j + \dots + w_k} \right). \quad (1.3.2)$$

### 1.3.1. Proposed Selection Procedure

Case I.  $\sigma_0^2$  known, common variance  $\sigma^2$  known, and common sample size  $n$ .

Definition 1.3.1. We define the procedure  $\pi_1$  as follows:

Step 1. Select  $\pi_i$ ,  $i = 1, \dots, k$  and stop, if

$$\hat{x}_{1:k} \geq \pi_0 - d_{1:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\pi_1$  and go to step 2.

Step 2. Select  $\pi_i$ ,  $i = 2, \dots, k$  and stop, if

$$\hat{x}_{2:k} \geq \pi_0 - d_{2:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\pi_2$  and go to step 3.

:

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Step  $k-1$ . Select  $\pi_i$ ,  $i = k-1, k$  and stop, if

$$\hat{x}_{k-1:k} \geq \pi_0 - d_{k-1:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\pi_{k-1}$  and go to step  $k$ .

Step  $k$ . Select  $\pi_k$  and stop, if

$$\hat{x}_{k:k} \geq \pi_0 - d_{k:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\pi_k$ .

Here  $d_{i:k}^{(1)}$ 's are the smallest values such that  $\pi_1 \in \Delta$ , that is,  $\pi_1$  satisfies the  $p^*$ -condition.

1.3.2. On the Evaluation of  $\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i-1})$  and the Value of the

Constants  $d_{1:k}^{(1)}, \dots, d_{k:k}^{(1)}$

For any  $\mu \in \mathcal{C}_i$ ,  $1 \leq i \leq k$ , let  $Z_i$ 's i.i.d.  $\mathcal{F}(\cdot; 0, 1)$  then

$$\begin{aligned} P_\mu(CS|_{i-1}) &= P_\mu\left(\bigcup_{j=1}^{k-i+1} \{X_{j:k} \leq \mu_0 + d_{j:k}^{(1)} \frac{\mu_j}{\sqrt{n}}\}\right) \\ &= P_\mu\left(\bigcup_{j=1}^{k-i+1} \bigcup_{r=1}^j \{X_{r:k} \leq \mu_0 + d_{j:k}^{(1)} \frac{\mu_r}{\sqrt{n}}\}\right) \\ &= P_\mu\left(\bigcup_{j=1}^{k-i+1} \bigcup_{r=1}^j \{Z_{r:k} \leq \frac{\mu_r - \mu_0}{\sqrt{n}} + d_{j:k}^{(1)}\}\right) \end{aligned}$$

which is decreasing in  $\mu_r$ ,  $r = 1, \dots, k-i+1$ .

Hence

$$\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i-1}) \geq P(Z_{k-i+1:k} \leq -d_{k-i+1:k}^{(1)})$$

On the other hand,

$$\begin{aligned} \inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i-1}) &= P_\mu\left(\bigcup_{j=1}^{k-i+1} \{X_{j:k} \leq \mu_0 + d_{j:k}^{(1)} \frac{\mu_j}{\sqrt{n}}\}\right) \\ &= P(Z_{k-i+1:k} \leq -d_{k-i+1:k}^{(1)}) \end{aligned}$$

whenever  $\mu^* = (\mu_0, -\frac{1}{\sqrt{n}}, \dots, -\frac{1}{\sqrt{n}}, \underbrace{\mu_0, \dots, \mu_0}_i)$

Thus, we have

$$\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i-1}) = P(Z_{k-i+1:k} \leq -d_{k-i+1:k}^{(1)}).$$

Since  $\hat{z}_{k-i+1:k} = \min \{z_{k-i+1}, \dots, z_{k-i+1+i} + \dots + z_k\}$  have the same distributions as

$$\hat{z}_{1:i} = \min \{z_1, \dots, z_{i-1} + z_i\},$$

let

$$\begin{aligned} V_i &= \hat{z}_{1:i} \\ &= \min_{1 \leq j \leq i} \frac{1}{r} \sum_{j=1}^r z_j, \end{aligned} \quad (1.3.4)$$

we have

$$\inf_{x \in \mathbb{R}} P_{\mu}(\text{CS}^i | s_1) = P(V_i \sim -d_{k-i+1:k}^{(1)}), \quad i = 1, \dots, k. \quad (1.3.4)$$

**Theorem 1.3.1.** In case I, ( $\omega_0$  known, common known  $\sigma^2$  and common sample size  $n$ ), if  $d_{k-i+1:k}^{(1)}$  is the solution of equation

$$P(V_i \sim -x) = P^*, \quad (1.3.5)$$

where

$$V_i = \min_{1 \leq j \leq i} \frac{1}{r} \sum_{j=1}^r z_j \quad \text{and} \quad z_i \text{ are i.i.d. } F(\cdot),$$

$i = 1, \dots, k$  then  $\hat{s}_1$  satisfies the  $P^*$ -condition.

**Proof.** For any  $i$ ,  $1 \leq i \leq k$ ,

$$\inf_{x \in \mathbb{R}} P_{\mu}(\text{CS}^i | s_1) = P(V_i \sim -d_{k-i+1:k}^{(1)}) = P^*,$$

so  $\hat{s}_1$  satisfies the  $P^*$ -condition.

Therefore, the problem of finding the  $d_{i:k}^{(1)}$ 's reduced to finding the distributions of  $V_1, \dots, V_k$ . This is achieved by using some theorems in the theory of random walk.

### 1.3.3. Some Theorems in the Theory of Random Walk

Suppose  $Y_1, Y_2, \dots$  are independent random variables with a common distribution  $B$  not concentrated on a half-axis, i.e.  $0 < P(Y_1 = 0) < 1$ ,  $P(Y_1 > 0) > 1$ . The induced random walk is the sequence of random variables

$$S_0 = 0, S_n = Y_1 + \dots + Y_n, \quad n = 1, 2, \dots$$

Let

$$\tau_n = P(S_1 \geq 0, \dots, S_{n-1} \geq 0, S_n \geq 0) \quad (1.3.6)$$

and

$$\tau(s) = \prod_{n=1}^{\infty} \tau_n s^n, \quad 0 \leq s \leq 1. \quad (1.3.7)$$

Then we have the following theorem which was discovered by Anderson (1953). Feller (1971) gave an elegant short proof.

Theorem 1.3.2.

$$\log \frac{1}{1-\tau(s)} = \sum_{n=1}^{\infty} \frac{s^n}{n} P(S_n = 0). \quad (1.3.8)$$

Theorem 1.3.3. (Feller (1971))

Let

$$p_n = P(S_1 = 0, \dots, S_n = 0),$$

then

$$p(s) = \prod_{n=1}^{\infty} p_n s^n = \frac{1}{1-\tau(s)}, \quad (1.3.9)$$

hence

$$\log p(s) = \sum_{n=1}^{\infty} \frac{s^n}{n} p(S_n = 0). \quad (1.3.10)$$

By symmetry, the probabilities

$$q_n = P(S_1 \leq 0, \dots, S_n \leq n) \quad (1.1.11)$$

have the generating function  $q$  given by

$$\log q(s) = \sum_{n=1}^{\infty} \frac{s^n}{n} P(S_n < 0). \quad (1.3.11)$$

Note: The above two theorems remain valid if the signs  $\wedge$  and  $\vee$  are replaced by  $\wedge$  and  $\vee$ , respectively.

Now, let

$$U_j^* = \max_{1 \leq i \leq j} \frac{1}{r} \sum_{i=1}^r z_i^*, \quad j = 1, 2, \dots, \quad (1.1.14)$$

and

$$V_j^* = \min_{1 \leq r \leq j} \frac{1}{r} \sum_{i=1}^r Z_i^*, \quad j = 1, 2, \dots, \quad (1.1.14)$$

where  $Z_i$ 's are i.i.d. with absolutely continuous c.d.f.  $g(\cdot)$ . We would like to apply Theorem 1.3.3 to get the distribution of  $U_i$  and  $V_i$ ,  $i = 1, 2, \dots$ .

Remark 1.3.1. The distribution of  $U_j^k$ ,  $j = 1, \dots, k$  for some  $k \leq 1$ , will be used whenever our goal is changed to select a subset containing no population with parameter smaller than the control.

**Theorem 1.3.4.** The generating function  $q(s)$  of  $P(I_i^t = x)$ ,  $i = 1, \dots$ , is

$$\lim_{n \rightarrow \infty} \frac{1}{n} \sum_{i=1}^n \mathbb{P}(X_i = 0) = \mathbb{P}(X_1 = 0) = 1 - \mathbb{E}X_1$$

where

$$S_n = \sum_{i=1}^n (Z'_i - x), \quad n = 1, 2, \dots,$$

if the distribution of  $Y_1 = Z'_1 - x$  is not concentrated on a half-axis.

Proof. Since the distribution of random variable  $Y_i = Z'_i - x$  is not concentrated on a half-axis, and  $Y_i$ 's are i.i.d. let  $S_r = \sum_{i=1}^r (Z'_i - x)$ ,  $r = 1, \dots, k$ . Then

$$(U_j > x) \geq \max_{1 \leq r \leq j} \frac{1}{r} S_r \geq 0^+ = \{S_1 \geq 0, \dots, S_j \geq 0\}.$$

By Feller's Theorem 1.3.3, we complete the proof.

Similarly,  $(V_j > x) = \{S_i \geq 0, \quad i = 1, 2, \dots, j\}$ ,

where

$$S_i = \sum_{r=1}^i (Z'_r - x).$$

Theorem 1.3.5. The generating function  $p(s)$  of  $P(V_j > x)$  is

$$\sum_{j=1}^{\infty} s^j P(V_j > x) = \exp \left[ \sum_{n=1}^{\infty} \frac{1}{n} s^n P(S_n > 0) \right], \quad (1.3.16)$$

if the distribution of  $Y_1 = Z'_1 - x$  is not concentrated on a half-axis.

Corollary 1.3.1. Both Theorem 1.3.4 and Theorem 1.3.5 hold for all  $x$  such that  $0 < G(x) < 1$ .

Proof. Let  $Y_1 = Z_1^t - x$ , then

$$P(Y_1 < 0) = G(x)$$

and

$$0 < G(x) < 1,$$

hence  $Y_1$  is not concentrated on a half-axis.

Corollary 1.3.2. Both Theorem 1.3.4 and Theorem 1.3.5 hold for all  $x$  whenever  $G = \phi$ , c.d.f. of  $N(0,1)$ , or  $G = \Gamma$  which is defined at the beginning of Section 1.3.

Proof. Followed immediately by Corollary 1.3.1.

Note that in the case of location parameter of normal population,

$$P(U_n^t < -x) = P(V_n^t > x).$$

Let

$$\gamma_j(x) - \gamma_j = P(S_j < 0), \quad j = 1, 2, \dots,$$

$$a(s) = \frac{\sum_{j=1}^n s^j}{n} \gamma_j,$$

we have

$$p(s) = \prod_{j=1}^n s^j P(V_j^t < x) = \exp(a(s)).$$

$$\text{Lemma 1.3.1. } p^{(n+1)}(s) = \sum_{j=0}^n \binom{n}{j} p^{(j)}(s) a^{(n+1-j)}(s), \quad n.$$

Proof. Since  $p'(s) = p(s) + a'(s)$ , the result can be proved by induction on  $n$ .

Theorem 1.3.6. Under the assumption of Theorem 1.3.5

$$\begin{aligned} P(V_{n+1}^+ < x) &= \frac{1}{(n+1)!} \lim_{s \rightarrow 0^+} \frac{d^{n+1}}{ds^{n+1}} p(s) \\ &= \frac{1}{n+1} \sum_{j=0}^n P(V_j^+ < x) \gamma_{n-j+1}, \quad n = 0, 1, 2, \dots \quad (1.3.17) \end{aligned}$$

where

$$P(V_0^+ < x) = 1, \quad \forall x.$$

Proof. By Lemma 1.3.1, we have

$$\begin{aligned} P(V_{n+1}^+ < x) &= \frac{1}{(n+1)!} \lim_{s \rightarrow 0^+} p^{(n+1)}(s) \\ &= \frac{1}{n+1} \sum_{j=0}^n \frac{1}{(n+1)!} \frac{n!}{j!(n-j)!} p^{(j)}(0) [(n-j)!] \gamma_{n+1-j} \\ &= \frac{1}{n+1} \sum_{j=0}^n \frac{p^{(j)}(0)}{j!} \gamma_{n+1-j} \\ &= \frac{1}{n+1} \sum_{j=0}^n P(V_j^+ < x) \gamma_{n+1-j}. \end{aligned}$$

Similarly, we have

$$P(U_{n+1}^+ < x) = \frac{1}{n+1} \sum_{i=0}^n P(U_{n-i+1}^+ < x) P(S_i = 0). \quad (1.3.18)$$

#### 1.3.4. Limiting Distributions of $U_n^+$ and $V_n^+$

Let  $f_n(x) = P(U_n^+ < x)$  and  $F_n(x)$  denote the limiting distribution function as  $n \rightarrow \infty$  of  $U_n^+$ . Suppose the distribution of random variable  $Y_1 - Y_1^+ < x$  is not concentrated on a half axis, then we have

$$1 - F_n(x) = P(S_1 > 0) + \sum_{r=2}^n P(S_1 > 0, \dots, S_{r-1} = 0, S_r = x),$$

$$1 - F_n(x) = \lim_{s \downarrow 1^-} \varphi(s),$$

and apply Andersen-Feller Theorem 1.3.2, we have

$$F_n(x) \approx \exp \left( - \sum_{r=1}^n \frac{1}{r} P(S_r > 0) \right). \quad (1.3.1)$$

Similarly,

$$G_n(x) \approx \exp \left( - \sum_{r=1}^n \frac{1}{r} P(S_r < 0) \right). \quad (1.3.2)$$

where

$$G_n(x) = P(V_n < x).$$

Let

$$G_n(-d_{1:n}^{(1)}) = P^*. \quad (1.3.3)$$

If  $Z_i$ ,  $i = 1, \dots, k$ , are independent identically distributed  $N(0,1)$ , then we can use the recurrence formula of Theorem 1.1.6 to solve the equations  $P(V_i < -d_{k-i+1:k}) = P^*$ ,  $i = 1, \dots, k$ . Hence in Case I,  $\varphi_j(x) = (-x)^j$ .

**Remark 1.3.2.** From formula (1.3.3) we know that  $d_{k-i+1:k}^{(1)}$  ( $i = 1, \dots, k$ ) does not depend on  $k$ . And we have  $d_{k-i+1:k}^{(1)} = d_{1:i}^{(1)}$ . These values for  $k = 1$  (1) 6, 10, ... and  $P^* = .99, .975, .95, .925, .9, .85, .75, .7, .65$  are tabulated in Table I.

### 1.3.5. Some Other Forms of Selection Procedure 1

Lemma 1.3.2.  $d_{1:i}^{(1)}$  is increasing in  $i$ .

Proof. By Remark 1.3.2 and the fact

$$v_{i+1} = \min (v_i, \frac{iv_i + z_{i+1}}{i+1}).$$

Lemma 1.3.3. If  $c_j$ ,  $1 \leq j \leq i \leq k$  is decreasing in  $j$ , then

$$\frac{\bigcup_{j=1}^i x_{j:k}}{i} - c_j = \frac{\bigcup_{j=1}^{i-1} x_{j:k}}{i-1} - c_j$$

Proof.  $\frac{\bigcup_{j=1}^i x_{j:k}}{i} - c_j = \frac{\bigcup_{j=1}^{i-1} x_{j:k}}{i-1} - c_j$ ,

since

$$x_{j:k} \leq \hat{x}_{j:k}, \quad 1 \leq j \leq k.$$

On the other hand, if

$$x_{r:k} = c_r \quad \text{for some } r, \quad 1 \leq r \leq i$$

then

$$x_{s:k} = c_r \quad \text{for some } s, \quad 1 \leq s \leq r,$$

since

$$x_{r:k} = \max_{1 \leq s \leq r} x_{s:k}.$$

Because  $c_j$  is decreasing in  $j$ , this implies  $x_{s:k} = c_r$  for some  $s$ ,  $1 \leq s \leq r$ .

Hence we have

$$\sum_{j=1}^i (x_{j:k} - c_j) = \sum_{j=1}^i (x_{j:k} - c_j'),$$

therefore the lemma is proved.

Definition 1.3.2. We define a selection procedure  $\cdot_1'$  by replacing the inequality in the  $i$ th step of procedure  $\cdot_1$  by the inequality

$$x_{i:k} \geq a_{i:k}^{(1)}, \quad i = 1, \dots, k$$

where  $a_{i:k}^{(1)}, \dots, a_{k:k}^{(1)}$  are the smallest values such that  $\cdot_1'$  satisfies the  $P^*$ -condition.

Theorem 1.3.7. The selection procedure  $\cdot_1$  and  $\cdot_1'$  are identical, i.e.  $a_{i:k}^{(1)} = a_{i:k}^1$ ,  $i = 1, \dots, k$ .

Proof. For any  $i$ ,  $1 \leq i \leq k$ , by Theorem 1.3.1

$$P^* = \inf_{\cdot \in \cdot_1} P_{\cdot}(\text{CS}^{(1)}_1) = P(Z_{k-i+1:k} \geq a_{k-i+1:k}^{(1)}).$$

On the other hand, using the same arguments as Section 1.3.7, we have

$$P^* = \inf_{\cdot \in \cdot_1'} P_{\cdot}(\text{CS}^{(1)}_1) = P(Z_{k-i+1:k} \geq a_{k-i+1:k}^{(1)}).$$

Hence we have  $a_{i:k}^{(1)} = a_{i:k}^1$ ,  $i = 1, \dots, k$ .

Since  $x_{1:k} = x_{i:k}$ , the first step of  $\cdot_1$  and  $\cdot_1'$  are identical. For  $i = 2, \dots, k$ , the event

$$\begin{aligned}
 \text{select } & \{i, \dots, k-1\} \cup \bigcup_{j=1}^i (x_{j:k} - 0 \cap d_{i:k}) \\
 & \cup \bigcup_{j=1}^i (x_{j:k} - 0 \cap d_{i:k}) \\
 & = \text{select } \{i, \dots, k-1\}
 \end{aligned}$$

by Lemma 1.3.2 and Lemma 1.3.3. Hence selection procedures  $\tau_1$  and  $\tau_2$  are identical.

### 1.3.6. Some Other Proposed Selection Procedures $\tau_2, \tau_3, \tau_4$

In Case I, we proposed some other selection procedures:

Definition 1.3.3. We define a selection procedure  $\tau_3$  by

$$\tau_3: \text{select } i \text{ if and only if } x_{i:k} - 0 \cap d_{i:k} \quad i = 1, \dots, k$$

where  $d$  is the smallest value such that  $\tau_3$  satisfies the  $\beta^*$ -condition.

Theorem 1.3.8. Under assumptions of Case I, and selection procedure  $\tau_3$ , if we select population  $i$ , then we will select populations  $j$ ,

for all  $j < i$ .

Proof. Since  $x_{i:k} \supset x_{j:k}$  for all  $j < i$ .

Evaluation the Value of  $\tau_3$

For any  $i, 1 \leq i \leq k$ , we have

$$\begin{aligned}
 \inf_{\tau_3} P(\text{CS}_{i-1}) & \leq \inf_{\tau_3} P(\text{CS}_{k-i+1:k}) \cap 0 \cap d_{i:k} \\
 & \leq P(V_i) = d^k
 \end{aligned}$$

by the same argument for selection procedure  $\gamma_1$  and here

$$V_i = Z_{1:i} - \min_{1 \leq j \leq i} \frac{1}{r} \sum_{j=1}^r Z_{j:i}.$$

We need the constant  $d$  such that  $P(V_i \geq -d) \geq P^*$  holds for all  $i$ ,

$1 \leq i \leq k$ . By Lemma 1.3.2 we have  $d = d_{1:k}^{(1)}$ . Hence we have the following theorem.

Theorem 1.3.3. Selection procedure  $\gamma_2$  satisfies the  $P^*$ -condition with  $d = d_{1:k}^{(1)}$ .

Corollary 1.3.3. If  $S_1$  and  $S_2$  are the selected subsets associated with selection procedures  $\gamma_1$  and  $\gamma_2$ , respectively, then  $\gamma_1 = \gamma_2$ .

Proof. Proof follows from Lemma 1.3.2.

Definition 1.3.4. The procedure  $\gamma_3$  is defined as follows:

Step 1. Select  $\pi_1$ ,  $i = 1$  and stop, if

$$X_1 - \mu_0 - d_1 \frac{\sigma}{\sqrt{n}} \geq 0,$$

otherwise reject  $\pi_1$  and go to step 2.

Step 2. Select  $\pi_2$ ,  $i = 2$  and stop, if

$$X_2 - \mu_0 - d_2 \frac{\sigma}{\sqrt{n}} \geq 0,$$

otherwise reject  $\pi_2$  and go to step 3.

⋮

Step  $k-1$ . Select  $\pi_k$ ,  $i = k-1$  and stop, if

$$X_{k-1} - \mu_0 - d_{k-1} \frac{\sigma}{\sqrt{n}} \geq 0,$$

otherwise reject  $x_{k+1}$  and go to step k.

Step k. Select  $x_k$  and stop, if

$$x_k - 0 \geq d_k \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $x_k$ .

Here  $x_j = \max(x_1, \dots, x_j)$  and  $d_j$ 's are the smallest values such that  $x_j$  satisfies the  $P^*$ -condition.

Evaluation of  $d_j$ 's

For any  $i$ ,  $1 \leq i \leq k$ ,

$$\begin{aligned} \inf_{u \in \mathcal{C}_i} P_u(CS^{1,3}) &= \inf_{u \in \mathcal{C}_i} P_u(x_{k+1} - x_i - u_0 \geq d_j \frac{\sigma}{\sqrt{n}}) \\ &= P_u(Z_{k-i+1} \geq d_{k-i+1} \frac{\sigma}{\sqrt{n}}) \\ &= P(Z_{k-i+1} \geq d_{k-i+1}) \\ &= P^*, \quad Z_i \sim F(\cdot; 0, 1). \end{aligned}$$

This implies  $d_{k-i+1} = d$  for all  $i$ , and

$$\begin{aligned} d &= F^{-1}(1-P^*), \\ &= F^{-1}(P^*), \quad \text{if } F \text{ is symmetric} \\ &= F^{-1}(P^*), \quad \text{if } X_i \sim N(x_i, \sigma^2/n). \end{aligned}$$

Similar to the selection procedure  $\gamma_1$ , we have the following theorem:

Theorem 1.3.10. Selection procedure  $\gamma_3$  satisfies the  $P^*$ -condition with

$$d_j = F^{-1}(1-P^*).$$

Definition 1.3.5. Selection procedure  $\tau_3^*$  is defined as follows:

Step 1. Select  $\tau_1$ ,  $i = 1$  and stop, if

$$x_1 - \mu \geq d \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\tau_1$  and go to step 2.

Step 2. Select  $\tau_2$ ,  $i = 2$  and stop, if

$$x_2 - \mu \geq d \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\tau_2$  and go to step 3.

⋮

⋮

Step  $k-1$ . Select  $\tau_k$ ,  $i = k-1$  and stop, if

$$x_{k-1} - \mu \geq d \frac{\sigma}{\sqrt{n}},$$

otherwise reject  $\tau_{k-1}$ .

Here

$$\begin{aligned} d &= F^{-1}(1-P^*) \\ &= F^{-1}(P^*) \text{ if } F \text{ is symmetric.} \end{aligned}$$

Theorem 1.3.11. The selection procedure  $\tau_3^*$  satisfies the P\* condition.

Proof. For any  $i$ ,  $1 \leq i \leq k$ ,

$$\inf_{\tau \in \tau_i} P_{\mu}(\tau S_i \geq 1) = P(\tau_{k-1} \geq 1) = P^*.$$

Theorem 1.3.12. The selection procedure  $\tau_3$  and  $\tau_3^*$  are identical.

Proof. The proof is simple hence it is omitted.

The following procedure  $\pi_4$  was given by Gupta and Sobel (1954), without assuming any ordering prior:

**Definition 1.3.6.** The selection procedure  $\pi_4$  is defined as follows:

$\pi_4$ : Select  $x_i$  if and only if  $x_i - d \geq 0$ ,  $i = 1, \dots, k$

where  $d$  is the smallest constant such that  $\pi_4$  satisfies the  $P^*$ -condition.

It was shown that the value  $d$  is determined by the equation

$$F(-d) = \frac{1}{1 + P^k}$$

or

$$F(d) = P^k \text{ if } F \text{ is symmetric.}$$

### 1.3.7. A Dual Problem

We start with the same assumptions as in Section 1.3.1 Case 1, but change our goal to select a subset which contains no bad populations; the definition of a correct selection (CS) will now be changed to select a subset that contains no bad populations and the  $P^*$ -condition will be defined based on this new definition of correct selection (CS).

In location parameter case, this problem is a dual problem of the original problem, namely, "select a subset which contains all good populations under ordering prior assumption".

One method to solve this problem is that, first, change the signs of all statistics and the control to opposite signs; then use a procedure for selecting a subset which contains all "new good" populations.

where the "new good" populations are the "old bad" populations, before changing signs; finally, reject the selected subset and keep the remainders as the desired selected subset. Let  $\omega_i$ ,  $i = 1, 2, \dots, 4$  denote the above procedure which corresponds to  $\omega_i$ ,  $i = 1, 2, \dots, 4$ , respectively.

**Theorem 1.3.3.** The selection procedure  $\omega_i$ ,  $i = 1, 2, \dots, 4$  satisfies the  $P^*$ -condition in which the correct selection (CS) means that it selects a subset which contains no bad population.

**Proof.** Given  $P^*$  and observations, for any selection procedure  $\omega_i$ ,  $i = 1, 2, \dots, 4$ , after changing the signs of all associated states, the probability that the selected subset  $S$  contains all "new good" populations is not less than  $P^*$ . If we reject the selected subset  $S$ , then the complement subset  $S^C$  of  $S$  contains any "new good" populations with probability less than  $1-P^*$ , but the "new good" populations are the originally bad populations so what we have is that the subset  $S^C$  contains any originally bad population with probability less than  $1-P^*$ , in other words, subset  $S^C$  contains no bad population with probability greater than or equal to  $P^*$ . Since this is true for all arbitrary true configurations, we have completed the proof.

**Remark 1.3.3.** It is easy to see that the value  $d_{i+1} \in \mathbb{N}$  of  $\omega_i$  which was used by  $\omega_i$  in the  $i$ th step is determined by the equation

$$P(\omega_{k-i+1} \mid d_{i+1} \in \mathbb{N}) = P^*,$$

where

$$\frac{P(\omega_{k-i+1} \mid d_{i+1} \in \mathbb{N})}{P(\omega_{k-i+1} \mid d_{i+1} \notin \mathbb{N})} = \frac{P^*}{1-P^*}.$$

If the distribution  $F$  is symmetric, then

$$d_{i;k}(.)_1 = d_{i;k}^{(1)}.$$

1.3.8. Some Proposed Selection Procedures  $\tau_i^{(2)}$ ,  $i = 1, 2, 3, 4$

When  $\theta_0$  is Unknown

Case 11.  $\theta_0$  unknown, common  $\sigma^2$  known, common sample size  $n$ .

Definition 1.3.7. We define a selection procedure  $\tau_1^{(2)}$  by requiring the inequalities

$$x_{i;k} - x_0 - d_{i;k}^{(1)} \geq 0, \quad i = 1, \dots, k$$

in procedure  $\tau_1$  (Definition 1.3.1) with

$$x_{i;k} - x_0 - d_{i;k}^{(2)} \geq 0, \quad i = 1, \dots, k, \text{ respectively.}$$

Here  $x_0 = \frac{n}{i+1} x_{0i}/n$ ,  $d_{i;k}^{(2)}$ ,  $i = 1, \dots, k$  are the smallest constants such

that the selection procedure  $\tau_1^{(2)}$  satisfies the  $P^*$ -condition.

Similar to the Case 1, we have the following theorem:

Theorem 1.3.14. For any  $i_1, \dots, i_k \in \{1, \dots, k\}$ ,  $\tau_{i_1, \dots, i_k}^{(2)}$  is determined by the equation

$$\int_{-\infty}^t F(V_{i_1} - t) + d_{k-i+1;k}^{(2)} dF(t) = P^*, \quad (1.3.14)$$

It is easy to see that  $d_{k-i+1:k}^{(2)} \geq d_{1:i}^{(2)}$ . The following theorem gives us an identical form of the selection procedure  $\psi_1^{(2)}$ .

**Theorem 1.3.15.** The selection procedure  $\psi_1^{(2)}$  will not be characterized if the statistics  $X_{i:k}$ ,  $i = 1, \dots, k$ , are replaced by  $\bar{X}_{i:k}$ ,  $i = 1, \dots, k$ , respectively.

**Proof.** The proof is the same as that in Case I and hence it is omitted.

The values  $d_{1:i}^{(2)}$ ,  $i = 1, \dots, k$  are tabulated in Table II for  $k = 1 (1) 6, 8, 10$ , and  $P^* = .99, .975, .95, .925, .90, .85, .75, .70, .65$ .

Similar to the Case I, we propose a selection procedure  $\psi_2^{(2)}$  as follows:

**Definition 1.3.8.** We define a selection procedure  $\psi_2^{(2)}$  by

$\psi_2^{(2)}$ : Select  $\tau_1$  if and only if  $\bar{X}_{i:k} - \bar{X}_0 = d_{1:i}^{(2)} \geq 1$ ,  $i = 1, \dots, k$

where  $d$  is the smallest value such that  $\psi_2^{(2)}$  satisfies the  $P^*$ -condition. Then, similar to Theorem 1.3.9 we have:

**Theorem 1.3.16.** Under assumptions of Case II, the selection procedure  $\psi_2^{(2)}$  satisfies the  $P^*$ -condition with  $d = d_{1:4}^{(2)}$ .

Next, we define a selection procedure  $\psi_3^{(2)}$  which is identical to  $\psi_2^{(2)}$  but replace  $\bar{X}_0$  by  $\tau_0$ , the sample mean of population  $\tau_0$ .

definition 1.3.9. The selection procedure  $\{x_i^{(2)}\}$  is defined by condition  $x_i = x_0 + d_i \frac{x_i - x_0}{\sqrt{n}}$  in  $\gamma_3$  (Definition 1.3.4) by  $x_i = x_0 + d_{i+1}^{(1)} \frac{x_i - x_0}{\sqrt{n}}$ ,  $i = 1, \dots, k$ , where  $d_1^{(1)}, \dots, d_k^{(1)}$  are the smallest values such that  $\{x_i^{(1)}\}$  satisfies the  $P^*$ -condition.

Similar to Theorem 1.3.10 we have:

Theorem 1.3.17. The selection procedure  $\{x_i^{(2)}\}$  satisfies the  $P^*$ -condition with  $d_i = d$ ,  $i = 1, \dots, k$  where  $d$  is determined by the equation

$$\int_{-\infty}^d [1 - F(t-d)]dF(t) = P^*, \quad (1.3.17)$$

$$\int_{-\infty}^d F(d-t)dF(t) = P^*, \text{ if } F \text{ is symmetric.}$$

And  $\{x_i^{(2)}\}$  will not be changed if the statistics  $x_i$  is replaced by  $\bar{x}_i$ , the sample mean of population  $\gamma_i$  for  $i = 1, \dots, k$ .

The following selection procedure  $\{x_i^{(3)}\}$  was proposed by Guntar and Sobel (1958):

Definition 1.3.10. The selection procedure  $\{x_i^{(3)}\}$  is defined by

$\{x_i^{(3)}\}$ : Select  $x_i^{(3)}$  if and only if  $x_i = x_0 + d \frac{x_i - x_0}{\sqrt{n_0}}$ ,  $i = 1, \dots, k$

where  $d$  is determined by the following equation if  $F$  is normal distribution:

$$\int_{-\infty}^d \int_{-\infty}^{x_i} [F(t) \int_{-\infty}^{x_0} F(u)du + d] F(u)du = P^*, \quad (1.3.18)$$

For the special case  $n_j = n$  ( $j = 0, 1, \dots, k$ )

$$\int_{-\infty}^{\infty} F^k(t+d)f(t)dt = P^k. \quad (1.3.25)$$

If  $F$  is normal distribution  $N(0,1)$ , the tables of  $d$ -values satisfying the Equation (1.3.25) for several values of  $P^k$  are given in Bechhofer (1954) for  $k = 1, 2, 3, 4$  and in Gupta (1946) for  $k = 1$ .

### 1.3.9. Some Proposed Selection Procedures $\{S_j^{(2)}\}$ , $j = 1, \dots, k$

When common Variance  $\sigma^2$  is unknown

Case III.  $\mu_0$  known, common variance  $\sigma^2$  unknown,  $n_j = n$ ,  $j = 1$

In this case, we assume that  $\sigma^2$  is  $\sigma^2(\sigma^2)$  which is the  $\sigma^2$  of  $N(0,1)$ .

**Definition 1.3.11.** We define the selection procedure  $\{S_j^{(3)}\}$  to replace the inequalities

$$c_{j(n)} = n + d_{j(n)}^{(3)} \quad j = 1, \dots, k$$

in procedure  $\{S_j\}$  (Definition 1.3.10) by

$$S_j^{(3)} = \left\{ \begin{array}{ll} S_j^{(2)} & \text{if } j = 1, \dots, k-1 \\ S_k^{(2)} & \text{if } j = k \end{array} \right. \text{respectively,}$$

where  $d^{(3)}$ 's are the smallest values such that  $S_j^{(3)}$  satisfies the  $P^k$ -condition,  $S_k^{(2)}$  denotes the selection function  $S_k$  for  $\sigma^2 = \sigma^2(n)$ , that is,

$$S_k = \frac{\sigma^2}{\sigma^2 + \frac{1}{n}} \quad \text{for } \sigma^2 = \sigma^2(n)$$

Note that  $\frac{\chi^2_{N-k}}{2}$  has the chi-square distribution  $\chi^2$  with  $N-k$  degrees of freedom.

By using similar arguments as in Case I, we have:

Theorem 1.3.18. The equation which determines the constant  $d_{k-i+1;k}^{(N)}$  is

$$P(V_i \leq d_{k-i+1;k}^{(N)}) = P^* \quad (1.3.18)$$

or

$$\int_0^{V_i} P(V_i \leq d_{k-i+1;k}^{(N)}) q_{(N)}(y) dy = P^* \quad (1.3.18)$$

where

$$V_i = \min_{1 \leq r \leq i} \frac{1}{r} \sum_{j=1}^r Z_j$$

and  $q_{(N)}(y)$  is the density of  $\frac{\chi^2_{N-k}}{2}$ .

We can rewrite formula (1.3.18) as

$$\int_0^{V_i} P(V_i \leq d_{k-i+1;k}^{(N)}) \sqrt{\frac{2}{\pi}} \int_0^{\infty} e^{-\frac{t}{2}} dt \frac{e^{-\frac{y^2}{t}}}{\sqrt{2\pi}} dt = P^* \quad (1.3.18)$$

or

$$\int_0^{V_i} P(V_i \leq d_{k-i+1;k}^{(N)}) \sqrt{\frac{2}{\pi}} \int_0^{\infty} e^{-\frac{t}{2}} dt \frac{e^{-\frac{y^2}{t}}}{\sqrt{2\pi}} dt = P^* \quad (1.3.18)$$

Remark 1.3.4. The values of  $d_{k-i+1;k}^{(3)}$ ,  $i = 1, \dots, k$  depend on  $\pi^{(3)}$ , hence  $d_{k-i+1;k}^{(3)} \neq d_{1;k}^{(3)}$ .

By using Rabinowitz and Weiss table (1959) (with  $n = 9$ ,  $k = 6$ ), we have evaluated and tabulated the values of  $d_{k-i+1;k}^{(3)}$ ,  $i = 1, \dots, k$ , in Table III, for  $k = 2$  (1) 6,  $\pi^* = .29, .35, .36, .37, .38, .39, .40$ , .45, .50, .60, and .75 with common sample size  $n = 3, 5, 9$ , and 21.

For  $k = 6$  and  $n = 21$ , we can use  $d_{1;k}^{(3)}$  as an approximation of  $d_{k-i+1;k}^{(3)}$ .

Definition 1.3.5. We define the selection procedure  $\hat{\gamma}_j^{(3)}$  by

$\hat{\gamma}_j^{(3)}$ : Select  $j$  if and only if  $\hat{\gamma}_{j;k}^{(3)} \geq d_{j;k}^{(3)}$ ,  $j = 1, \dots, n$ ,  
where  $\hat{\gamma}$  is defined as in procedure  $\hat{\gamma}_j^{(3)}$ , and  $d_{j;k}^{(3)}$  is the smallest constant such that  $\hat{\gamma}_j^{(3)}$  satisfies the  $P^*$ -condition.

As before, it can be shown that  $\hat{\gamma}_j^{(3)} = \hat{\gamma}_{j;k}^{(3)}$ .

Remark 1.3.5. Theorem 1.3.2 still holds for case III, i.e. the selection procedure  $\hat{\gamma}_j^{(3)}$  will not be changed if we replace the factor statistics  $\hat{\gamma}_{j;k}$  by  $\hat{\gamma}_{j;k}$ , respectively. But this is not necessarily true for selection procedure  $\hat{\gamma}_j^{(3)}$ .

Definition 1.3.6. The selection procedure  $\hat{\gamma}_j^{(3)}$  is defined to have the same form as procedure  $\hat{\gamma}_j^{(3)}$  except that the inequality defining the  $j$ th step of procedure  $\hat{\gamma}_j^{(3)}$  is replaced by

$$\hat{\gamma}_{j-1}^{(3)} \geq \alpha_j = \frac{1}{\sqrt{j}} \quad \text{for } j = 1, \dots, k.$$

The proof of the following theorem uses the same arguments as that in Case I, hence it is omitted.

Theorem 1.3.19. The equation which determines the constant  $d$  of selection procedure  $\tau_3^{(3)}$  is

$$\int_0^{\infty} \tau_3^{(3)}(yd) q_1(y) dy = P^*, \quad (1.3.31)$$

Gupta and Sobel (1958) gave a selection procedure  $\tau_4^{(4)}$  in this case. It is as follows:

$\tau_4^{(4)}$ : Select  $x_i$  if and only if  $x_i - x_0 + d \geq \frac{S}{m_i}$  for  $i = 1, \dots, k$

and the equation which determines  $d$  is

$$\int_0^{\infty} \tau_4^{(4)}(yd) q_1(y) dy = P^*. \quad (1.3.32)$$

1.3.10. Some Proposed Selection Procedures  $\tau_j^{(4)}$ ,  $j = 1, 2, 3, 4$

When Both Control  $\tau_0$  and Common Variance  $\sigma^2$  are Unknown.

Case IV.  $\tau_0$  unknown, common variance  $\sigma^2$  unknown and common sample size  $n$ .

We assume that in this case distribution  $F$  is the c.d.f.  $N(t, 1)$ , and denoted by  $\tau_1$ . We replace  $\tau_0$  in each selection procedure  $\tau_j^{(3)}$  by  $\tau_0$ ,  $1 \leq j \leq 3$ , and get four procedures  $\tau_j^{(4)}$ ,  $1 \leq j \leq 4$ , respectively. Let  $\chi^2(t)$  denote the c.d.f. of the chi-square distribution with  $k(n-1)$  degrees of freedom. The constant  $d_{k-i+1;k}^{(4)}$ ,  $i = 1, \dots, k$ , of procedure  $\tau_1^{(4)}$  is determined by

$$\int_0^1 \int_0^1 P(Y_1 = u + d_{k-i+1, k-i}^{(4)}(t)u, Y_2 = t) d u d t, \quad i = 1, \dots, k.$$

The constant  $d$  of procedure  $\psi_2^{(4)}$  is

$$d = d_{1, k}^{(4)}.$$

The constants  $d$  of procedures  $\psi_3^{(4)}$  and  $\psi_4^{(4)}$  are determined by

$$\int_0^1 \int_0^1 r(u + td) c^*(u) d u d t = P^*, \quad i = 3, 4,$$

with  $r = 1$  and  $k$ , respectively, and their values for selected values of  $P^*$ ,  $k$  and  $i$  are given in Gupta and Sebel (1967) and Bennett (1971).

### 1.3.11. Properties of the Selection Procedures

Under simple ordering prior, it is natural to require that an order selection procedure is order-preserving as defined below:

**Definition 1.3.14.** A selection procedure  $\psi$  is order-preserving if it selects  $\gamma_i$  with parameter  $\alpha_i$ , and if  $\gamma_i < \gamma_j$ , then it also selects  $\gamma_j$ . Procedure  $\psi$  is weak order-preserving or monotone if

$$P(\gamma_i \text{ is selected}) \geq P(\gamma_j \text{ is selected}) \text{ whenever } \gamma_i < \gamma_j.$$

It is easy to see that any order-preserving selection procedure is weak order-preserving, but the converse is not true.

Now, let  $\psi_i^{(1)} = \gamma_i$ ,  $i = 1, \dots, k$ .

**Theorem 1.3.20.** The selection procedures  $\psi_1^{(1)}$ ,  $\psi_2^{(1)}$  and  $\psi_3^{(1)}$  are order-preserving and procedure  $\psi_4^{(1)}$  is monotone for  $i = 1, \dots, k$ .

Proof. The proof follows immediately from the definition of the procedures.

Given observations  $X = x = (x_0, \dots, x_k)$  where  $x_i$  is the sample mean of population  $\gamma_i$ ,  $i = 1, \dots, k$ , and  $x_0 = \bar{x}_0$  if  $\gamma_0$  is known, otherwise  $x_0$  is the sample mean of population  $\gamma_0$ . Let

$$\gamma_i(x, \cdot) = P(\gamma_i \text{ included in the selected subset } x = x_1, \dots, x_k) \text{ for } i = 1, \dots, k.$$

**Definition 1.3.16.** A selection procedure  $\gamma$  is called translation-invariant if for any  $c \in \mathbb{R}^{k+1}$ ,  $c \in \mathbb{R}$

$$\gamma(x_0 + c, x_1 + c, \dots, x_k + c, \cdot) = \gamma(x_0, \dots, x_k, \cdot) \text{ for } i = 1, \dots, k.$$

**Theorem 1.4.21.** The selection procedures  $\gamma_1^{(i)}, \gamma_2^{(i)}, \gamma_3^{(i)}$  and  $\gamma_4^{(i)}$  are translation-invariant for  $i = 1, 2, 3, 4$ .

**Proof.** By Corollary 1.3.1 the isotonic regression is a linear operator. On the other hand,

$$\frac{\sum_{i=1}^n \gamma_i(x_i + c)}{n} = \frac{\sum_{i=1}^n \gamma_i(x_i)}{n} + c,$$

hence we have the result.

#### Expected Number (Size) of Bad Populations in the Selected Subsets

Suppose the control  $\gamma_0$  is known and we have common sample size  $n$  and common known variance  $\sigma^2$  without loss of generality we assume that  $\gamma_0 = 0$  and  $\gamma_0 = 1$ . Let  $\gamma(x)$  denote the expected number of bad populations in the selected subset  $x$  using the selection procedure

, then for any  $i$ ,  $0 \leq i \leq k$ ,

$$\begin{aligned} \sup_{\mu \in \mathcal{K}_{k-i}} E(S^{(1)}_{\mu, 1}) &= \sup_{\mu \in \mathcal{K}_{k-i}} \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\mu, j} \cap d_{1;k}^{(1)}) \\ &= \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\mu, j} \cap d_{1;k}^{(1)}) \quad (\text{from (1.3.35)}) \\ &= \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\mu, j} \cap d_{1;k}^{(1)}) \quad (\text{from (1.3.36)}) \end{aligned}$$

On the other hand, for procedure  $\varphi$ ,

$$\sup_{\mu \in \mathcal{K}_{k-i}} E(S^{(1)}_{\mu, 2}) = \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\mu, j} \cap d_{1;k}^{(1)}) \quad (\text{from (1.3.36)})$$

Formula (1.3.36) is increasing in  $i$  and is greater than or equal to

Formula (1.3.35), since

$$d_{1;k}^{(1)} \geq d_{1;k+1}^{(1)} \geq d_{1;k}^{(1)}.$$

Therefore, we have the following theorem.

Theorem 1.3.22. For any  $i$ ,  $0 \leq i \leq k$

$$\sup_{\mu \in \mathcal{K}_i} E(S^{(1)}_{\mu, 2}) \geq \sup_{\mu \in \mathcal{K}_i} E(S^{(1)}_{\mu, 1}),$$

$$\sup_{\mu \in \mathcal{K}_0} E(S^{(1)}_{\mu, 2}) \geq \sup_{\mu \in \mathcal{K}_0} E(S^{(1)}_{\mu, 1}).$$

Theorem 1.3.23. In Section 1.3.1, Case 1, for any  $i$ ,  $0 \leq i \leq k$

$$\sup_{\mu \in \mathcal{K}_{k-i}} E(S^{(1)}_{\mu, 2}) = \gamma + \eta(1-\eta^i)/P^* \quad (\text{from (1.3.36)})$$

where  $\eta = 1 - P^*$ .

Proof.

$$\begin{aligned}
 & \sup_{\mathcal{E} \subset [k-j]} E(S^{t+1}_{\mathcal{E}}) \\
 &= \sup_{\mathcal{E} \subset [k-j]} \sum_{i=1}^j P(\text{select } r_i^{t+1} \text{ in } \mathcal{E}) \\
 &= \sup_{\mathcal{E} \subset [k-j]} \sum_{i=1}^j P(\max_{1 \leq r \leq i} r = d) \\
 &= \sum_{i=1}^j (1 - \frac{i}{r+1}) P(-d) \\
 &= j - \sum_{i=1}^j q^i \\
 &= j - q(1-q^j)/P^*.
 \end{aligned}$$

where  $q = (1-P^*)$ .

Theorem 1.3.24.  $\sup_{\mathcal{E} \subset [k-j]} E(S^{t+1}_{\mathcal{E}})$  is increasing in  $j$ , hence

$$\sup_{\mathcal{E} \subset [k-j]} E(S^{t+1}_{\mathcal{E}}) \geq \sup_{\mathcal{E} = \emptyset} E(S^{t+1}_{\mathcal{E}}) = k - q(1-q^k)/P^*. \quad (1.3.25)$$

Proof. Since the function

$$f(x) = x - ab^{x+1}$$

is increasing in  $x$ , for  $0 < a < 1$ ,  $0 < b < 1$ , and  $0 < x < \infty$ .

In Case I of Gupta (1965) showed that

$$\sup_{\mathcal{E}} E(S^{t+1}_{\mathcal{E}}) = kp^*k. \quad (1.3.26)$$

When the ordering prior about the unknowns is either unknown or unknown, we can use the selection procedure of Gustafsson and Broström (1977) using the ordering of the sample means as the ordering of unknowns, or we can and apply the selection procedure which is currently used in the literature prior. With the latter approach, the substitution is made that the isotonic regression of the sample means starts to the left of the true sample means, and that the selection procedures  $\pi_{ij}^{(1)}, i = 1, 2, 3, j = 1, 2, 3$  and  $\pi_{ij}^{(2)}, i = 1, 2, 3, j = 2, 3$  are the same as  $\pi_{ij}^{(3)}, i = 1, 2, 3, j = 4$ , respectively, and the selection procedures  $\pi_{ij}^{(1)}, i = 1, 2, 3, 4$  for  $\pi_{ij}^{(4)}, i = 1, 2, 3, 4$  are the same, respectively, which are equivalent to the procedures proposed by Gustafsson (1975) and Broström (1977), independent 1.

It has been proved in some quite general situation, and it has been proved by using Monte Carlo technique in some selected cases by Gustafsson (1975) and Broström (1977), separately, that  $\pi_{ij}^{(1)} - \pi_{ij}^{(3)}$  is slightly better than  $\pi_{ij}^{(4)}$ . The values  $d_{ij,k}^{(j)}$  in the  $j$ th step of the procedure  $\pi_{ij}^{(1)}$ ,  $i = 1, 2, 3, j = 1, 2, 3, 4$ , are given by Broström (1977) as follows:

$$d_{ij,k}^{(1)} = \frac{1}{\pi_{ij}^{(1)}(1 - (\pi_{ij}^{(1)})^k)}, \quad (1.1)$$

$$\int_{-\infty}^{x_i} \pi_{ij}^{(1)}(x + d_{ij,k}^{(1)}) \phi(x) dx = \pi_{ij}^{(1)}, \quad (1.2)$$

$$\int_0^{\infty} \pi_{ij}^{(1)}(x d_{ij,k}^{(1)}) \phi(x) dx = \pi_{ij}^{(1)}, \quad (1.3)$$

and

$$\int_{0}^{x_i} \int_{-\infty}^{x_i} \pi_{ij}^{(1)}(x + y d_{ij,k}^{(1)}) \phi(x) \phi(y) dy dx = \pi_{ij}^{(1)}, \quad (1.4)$$

where  $\phi(\cdot)$  is the density of  $N(0, 1)$ .

### 3.1. Selection Procedure for Scale Parameter of Gamma Population

Suppose we have  $k+1$  independent populations  $\pi_1, \pi_2, \dots, \pi_{k+1}$ . The population  $\pi_i$  has a gamma density function

$$g(x_i | \alpha_i, \beta_i) = \frac{1}{\beta_i^{\alpha_i}} \frac{x_i^{\alpha_i-1}}{\Gamma(\alpha_i)} e^{-x_i/\beta_i}, \quad x_i > 0.$$

Example. The ordering prior of  $\alpha_1, \alpha_2, \dots, \alpha_{k+1}$  is  $\alpha_1 \leq \alpha_2 \leq \dots \leq \alpha_{k+1}$ , i.e.,  $\alpha_1 = 0, \alpha_2 = 1, \dots, \alpha_{k+1} = p$ . Note that the values of  $\beta_i$ 's and  $\alpha_i$ 's are unknown,  $\alpha_i$ 's are known.

In this section we define population  $\pi_i$  to be the  $i$ -th most scale population if the scale parameter  $\beta_i = \beta_{\pi_i}$ . Let  $\alpha_i$  be  $\alpha_i$ 's, then the parameter space is denoted by  $\mathbb{R}^k$ , where

$$\{\alpha \in \mathbb{R}^{k+1} : \alpha_1 < \alpha_2 < \dots < \alpha_{k+1}\}$$

is a subspace of  $(k+1)$ -dimension Euclidean space  $\mathbb{R}^{k+1}$ .

Suppose we have independent observations  $x_{ij}$  ( $i = 1, \dots, n_j$ ) from population  $\pi_j$  ( $i = 1, \dots, k$ ). Let  $\bar{x}_{ij} = \bar{x}_{ij} / n_j$ , then

$$x_{ij} - \frac{n_j}{j+1} \bar{x}_{ij} / n_j \text{ has density } g(x_{ij} | \alpha_j, \beta_{\pi_j}),$$

and

$$\bar{x}_{ij} / n_j \text{ has density } g(\bar{x}_{ij} | \alpha_j, \beta_{\pi_j}).$$

Suppose our goal is to select a subset which contains all the populations under the ordering prior with probability greater than or equal to  $P_k$ , a predecision value between zero and one.

Let  $\pi_{\pi_1} \cup \dots \cup \pi_{\pi_k}$  be the subspace of parameter space  $\mathbb{R}^k$  such that  $\pi_{\pi_i} = \{ \alpha \in \mathbb{R}^k : \alpha_i = \pi_i \}$  where

$$\begin{aligned}
 \hat{\gamma}_i &= \min(\hat{\gamma}_i, \hat{\gamma}_0 \gamma_{i+1}), \quad \text{if } i = 1, \dots, m-1, \\
 \hat{\gamma}_k &= \hat{\gamma}_k \gamma_0, \quad \text{if } i = k, \\
 \hat{\gamma}_0 &= \hat{\gamma}_0 \gamma_1, \quad \text{if } i = 0.
 \end{aligned}$$

#### 1.4.1. Proposed Selection Procedures $\hat{\gamma}_i$ , $i = 6, 7, 8, 9$

Case I. Control  $\gamma_0$  known and common sample size  $n$ .

Definition 1.4.1. The selection procedure  $\hat{\gamma}_6$  is defined as follow

Step 1. Select  $\hat{\gamma}_i$ ,  $i = k$  and stop, if

$$\hat{\gamma}_{k:k} \leq c_{k:k} \gamma_0$$

otherwise reject  $\hat{\gamma}_k$  and go to step 2.

Step 2. Select  $\hat{\gamma}_i$ ,  $i = k-1$  and stop, if

$$\hat{\gamma}_{k-1:k} \leq c_{k-1:k} \gamma_0$$

otherwise reject  $\hat{\gamma}_{k-1}$  and go to step 3.

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Step k-1. Select  $\hat{\gamma}_i$ ,  $i = 2$  and stop, if

$$\hat{\gamma}_{2:k} \leq c_{2:k} \gamma_0$$

otherwise reject  $\hat{\gamma}_2$  and go to step k.

Step k. Select  $\hat{\gamma}_1$  and stop, if

$$\hat{\gamma}_{1:k} \leq c_{1:k} \gamma_0$$

otherwise reject  $\hat{\gamma}_1$ .

where  $\epsilon_{ijk}(-1)$ ,  $i = 1, \dots, s$ , are the smallest values such that the procedure  $\rho$  satisfies the  $\mathcal{P}^*$ -condition.

**Theorem 1.4.1.** Assume we have a uniform sample size  $n$  and  $\epsilon_1, \epsilon_2, \dots, \epsilon_n \in \{-1, 1\}$ , and the constant  $c_{\text{opt}} \in [0, 1]$  is determined by the condition

$$P(\Omega_j = e_{i,j,k}) = P^*, \quad j = 1, \dots, k \quad (1.10)$$

Wittman

$$U_i = \max_{1 \leq s \leq i} \left( \frac{Y_s + \dots + Y_i}{i-s+1} \right) \quad (1.1)$$

and  $Y_i$  are i.i.d. with density

and  $\text{cal. f. } \mathcal{U}(\cdot, \cdot, \cdot, \cdot)$

then the procedure  $\mathcal{G}$  satisfies the  $\mathcal{B}^*$ -condition.

Proof. For any  $i$ ,  $1 \leq i \leq k$ , if the unknown true value  $\theta_i$  is fixed, then there are  $\lceil k \rceil$  good populations, then, under the procedure  $\tau_i$ ,

$$\begin{aligned}
 & \inf_{\substack{c \in \mathbb{R} \\ c \geq 0}} \mathbb{P} \left( \mathcal{C}^{n+1}_0 \in \cdot \right) \\
 &= \inf_{\substack{c \in \mathbb{R} \\ c \geq 0}} \mathbb{P} \left( \mathcal{C}^{n+1}_0 \in \cdot \mid \mathcal{C}^n_0 = \frac{1}{2} \right) \\
 &= \inf_{\substack{c \in \mathbb{R} \\ c \geq 0}} \mathbb{P} \left( \mathcal{C}^{n+1}_0 \in \cdot \mid \mathcal{C}^n_0 = \frac{1}{2} \text{ and } \mathcal{C}^n_1 = \frac{1}{2} \right) \\
 &= \inf_{\substack{c \in \mathbb{R} \\ c \geq 0}} \mathbb{P} \left( \mathcal{C}^{n+1}_0 \in \cdot \mid \mathcal{C}^n_0 = \frac{1}{2} \text{ and } \mathcal{C}^n_1 = \frac{1}{2} \text{ and } \mathcal{C}^n_2 = \frac{1}{2} \right) \\
 &= \dots
 \end{aligned}$$

$$= \inf_{\mathbf{c}_{i:k}} P_{\mu} \left( \bigcup_{j=i}^k \left( \max_{1 \leq s \leq j} \min_{s \leq t \leq k} \frac{Y_{s-1} + \dots + Y_{t-1}}{t-s+1} \leq c_{i:k} \right) \right)$$

$$= P_{\mu \otimes \nu} \left( \bigcup_{j=i}^k \left( \max_{1 \leq s \leq j} \min_{s \leq t \leq i} \frac{Y_{s-1} + \dots + Y_{t-1}}{t-s+1} \leq c_{i:k} \right) \right)$$

$$= P \left( \max_{1 \leq s \leq i} \frac{Y_{s-1} + \dots + Y_i}{i-s+1} \leq c_{i:k} \right)$$

$$= P(U_i \leq c_{i:k}),$$

where  $Y_i$ 's are i.i.d with density  $a(\cdot; \nu, \frac{1}{n})$ ,

$$\mathbf{v}^* = (\underbrace{0, 0, \dots, 0}_{i+1}, \dots, \dots)$$

and

$$U_i = \max_{1 \leq s \leq i} \frac{Y_{s-1} + \dots + Y_i}{i-s+1}.$$

**Corollary 1.4.1.**  $c_{i:k} = c_{i:i}$ ,  $i = 1, \dots, k$ .

For any  $x > 0$ , let  $S_n = \sum_{i=1}^n (Y_i - x)$ ,  $n = 1, 2, \dots$ ,  $c_{ij} = 0$  if  $i > j$ .

$0 < P(Y_i - x > 0) < a(x, \nu, \frac{1}{n}) < 1$ , the distribution of  $Y_i - x$  is concentrated on a half-axis. By Theorem 1.4.1, the probability generating function of cumulative distribution functions  $P(U_i \leq \cdot)$ ,  $i = 1, \dots, k$ , is given by

$$\exp \left[ \frac{1}{\sqrt{1-\nu}} \left( \frac{1}{k} + \frac{1}{n} \right) + a(kn, \nu, \frac{1}{n}) \right].$$

Hence by Theorem 1.3.6, we have the following recurrence formula for the all  $x > 0$

$$P(U_{k+1} = x)$$

$$= \sum_{j=0}^{k+1} \sum_{i=0}^j P(U_{k+1} = x) + G((i+1)n; (j+1) + 1) \quad (1.4.4)$$

where

$$P(U_0 = x) = 1.$$

When  $x = 0$ , both sides of Equation (1.4.4) equal to zero, hence it also holds for  $x = 0$ .

Note that

$$\begin{aligned} P\left(\frac{1}{r} \sum_{i=1}^r Y_i = x\right) &= G(x; r, \frac{1}{rn}) \\ &= G(xn; r, 1). \end{aligned} \quad (1.4.5)$$

The values  $c_{k:k}(P^*, r, n)$  which satisfy Equation (1.4.1) are tabulated in Table IV for  $k = 1$  (1) 10,  $P^* = .29, .45, .60, .75,$   $.82, .4, .6$ , and  $n = 4, 6, 8, 10, 15, 20$ .

Lemma 1.4.4.  $c_{i:k} = c_{i+1:k}$  for all  $1 \leq i \leq k-1$ .

Proof. The constants  $c_{i:k}$  ( $i = 1, \dots, k$ ) are determined by (1.4.1), respectively.

$U_{i+1} = U_i + S_i$  implies  $c_{i:k} = c_{i+1:k}$  for all  $1 \leq i \leq k-1$ .

Theorem 1.4.7. The selection procedure  $\varphi$  will not be changed if the monotone estimators  $\hat{Y}_{i:k}$ ,  $i = 1, \dots, k$  are replaced by  $\hat{c}_{i:k} \hat{Y}_{i+1:k}$ ,  $i = 1, \dots, k-1$ .

where

$$\begin{aligned} Y_{j+1} &= \max_{1 \leq i \leq j} Y_{i+1} \\ &= \max_{1 \leq i \leq j} \frac{f_{S_i} + \epsilon_i}{1 - \epsilon_i} \end{aligned}$$

Proof. The proof is similar to that of Theorem 1.4.1.

Next, we define a selection procedure by using an efficient estimator and a fixed constant which depends on  $P^*$ ,  $F_0$ , sample size  $n$  and constant  $c_3 < 1$ .

**Definition 1.4.2.** The selection procedure  $\gamma_j$  is defined by

$\gamma_j$ : Select  $x_j$  if and only if  $Y_{j+1} < c_3$  for  $j = 1, \dots, n$ ,

where  $c_3 \in (1)$  is the smallest value such that procedure  $\gamma_j$  satisfies the  $P^*$ -condition.

**Corollary 1.4.2.** The constant  $c_3(c(P^*, F_0))$  of the selection procedure  $\gamma_j$  equals to  $c_{k;k}$  which is determined by the following

Proof. Follows immediately from Theorem 1.4.1 and (1.4.1).

**Definition 1.4.3.** The selection procedure  $\gamma$  is defined as follows:

Step 1: Select  $x_1$  if and only if  $Y_2 < c_3$ .

$$\begin{aligned} Y_3 &= \max_{1 \leq i \leq 2} Y_{i+1} \\ &= \max_{1 \leq i \leq 2} \frac{f_{S_i} + \epsilon_i}{1 - \epsilon_i} \end{aligned}$$

otherwise proceed to Step 2.

Step 2. Select  $c_j$ ,  $j = k+1$ , and stop, if

$$\begin{cases} x_{k+1} < c_{k+1} \\ c_{k+1} < 0 \end{cases}$$

otherwise, reject  $c_{k+1}$  and go to step 1.

Step 3.1. Select  $c_j$ ,  $j = 1$ , and stop, if

$$\begin{cases} x_1 < c_1 \\ c_1 < 0 \end{cases}$$

otherwise, reject  $c_1$  and go to step k.

Step k. Select  $c_1$  and stop, if

$$\begin{cases} c_1 < 0 \\ c_1 < 1 - \beta \end{cases}$$

otherwise, reject  $c_1$ .

where the  $c_j$ 's are the smallest real values (if any) such that the  $(x_j, c_j)$  vector satisfies the  $\beta^*$ -constraint.

where  $\beta^* \in [0, 1]$ , the constraint  $c_j$  are determined by

$$\begin{cases} 1 - \beta^* \leq x_j < c_j < 1 - \beta, & \text{if } j = 1, \dots, k-1 \\ c_j < 0, & \text{if } j = k \end{cases}$$

and the optimality is checked by  $\delta_j < \delta_{j+1}$ .

where

$$\delta_j =$$

$$= \inf_{\mathbf{c} \in \mathcal{C}} P_{\mathbf{c}}(\bigcup_{i=1}^k (z_i \in [c_i, c_{i+1})))$$

$$= \inf_{\mathbf{c} \in \mathcal{C}} P_{\mathbf{c}}(\bigcup_{i=1}^k (z_i \in [c_i, c_{i+1})))$$

$$= P_{\mathbf{c}^*}(\bigcup_{i=1}^k (z_i \in [c_i, c_{i+1}]))$$

$$= P(z_i \in [c_i, c_{i+1}])$$

where  $z_i \sim \frac{e^{\lambda_i}}{\lambda_i} \mathcal{E}_i$ ,  $i = 1, \dots, k$  are i.i.d. with the same density.

$g(\cdot; \mathbf{c}_i, 1), \mathbf{c}^* = (\underbrace{c_0, \dots, c_0}_{i+1}, \dots, c_k)$ . Hence  $c_j$ ,  $j = 1, \dots, k$  are determined by (1.4.6). If  $c_1 < \dots < c_k$ , then  $c_1 = \dots = c_k$ .

The following selection procedure  $\pi_0$  was given by Gupta and Edel (1958).

**Definition 1.4.4.** The selection procedure  $\pi_0$  is defined by

(a) Select  $\pi_i$  if and only if  $\frac{c_i}{\lambda_i} \geq \frac{c_j}{\lambda_j}$ ,  $i = 1, \dots, k$   
where  $c^*$  is determined by

$$\frac{1}{\lambda_1} \frac{1}{\lambda_2} \dots \frac{1}{\lambda_k} \frac{c^*}{\lambda_1} \dots \frac{c^*}{\lambda_k} \in \mathcal{C} = \{ \mathbf{c} \in \mathbb{R}^k : c_i \geq 0, i = 1, \dots, k \}$$

for  $\pi_i \in \{1, \dots, k\}$ , it turns to

$$\frac{1}{\lambda_1} \frac{1}{\lambda_2} \dots \frac{1}{\lambda_k} \frac{c^*}{\lambda_1} \dots \frac{c^*}{\lambda_k} \in \mathcal{C} = \{ \mathbf{c} \in \mathbb{R}^k : c_i \geq 0, i = 1, \dots, k \}$$

the left-hand side is the c.d.f. of  $\frac{1}{2} \chi^2$  with  $r$  degrees of freedom, hence the value  $c^*$  can be easily solved with the help of a table of chi-square distribution.

#### Application to the Selection of Variance of Normal Population

$\Omega_{ij} = \sigma_{ij}^2$ ,  $i = 0, 1, \dots, k$  are the scale parameters for the  $k+1$  normal populations and  $y_{ij}$  ( $j = 1, \dots, n$ ;  $i = 1, \dots, k$ ) are the observations on the population  $\gamma_i$  with the mean  $\gamma_0$  (known). We assume that the order  $\sigma_1^2 < \dots < \sigma_k^2$  is known.

In the application of selection procedure  $\epsilon$  or  $\gamma_2$ , what we need to do is to evaluate the isotonic regression of  $S_i^2$  which is the sample variance of population  $\gamma_i$ ,  $i = 1, \dots, k$  and denote it by  $S_{i+k}^2$ ,  $i = 1, \dots, k$ , then directly apply  $\epsilon_6$  or  $\gamma_2$ . The constant we need is determined by Equations (1.4.2) and (1.4.4) where we replace  $n$  by  $n$ , the reason being that  $(nS_i^2/\sigma_i^2)$  has  $\chi^2_n$  distribution with  $n$  degrees of freedom and  $\epsilon_{ij} = S_{i+k}^2/\sigma_i^2$  has the c.d.f.

$$G(2nt; n, \beta) = G(t; n, \frac{1}{n}),$$

hence

$$P\left(\frac{1}{n} \sum_{i=1}^r y_{ij} - t\right) = G(tn; rn, 1).$$

The application of  $\gamma_2$  is similar to that of  $\epsilon_6$  (see Indiana Statist. (1959)). What we need to do is to replace  $y_{ij}$  in  $\gamma_2$  and  $\epsilon_{ij}$  by  $y_{ij} - t$ , replace  $n$  in Equation (1.4.6) and (1.4.7) by  $n$ ,  $i = 1, \dots, k$ .

Remark 1.4.1.  $\gamma_0$  (Gupta and Sobel (1958)) does not depend on the underlying prior and the sample sizes for each population need not be equal.

If the means  $\mu_i$ ,  $i = 1, \dots, k$  are unknown and common, and the sample size  $n = 1$ , let  $S_i^{(2)} = \sum_{j=1}^n (X_{ij} - \bar{X}_i)^2/n-1$  and use  $n = 1$  in place of  $n$  in (1.4.4), (1.4.6) and (1.4.8) which determined the constants  $c_{i;k}$ ,  $c$  and  $c'$  for  $\gamma_6$ ,  $\gamma_8$  and  $\gamma_9$ , respectively.

#### 1.4.2. Selection Procedure $\gamma_i^{(2)}$ , $i = 6, 7, 8, 9$

Case II.  $\gamma_0$  unknown.

The assumptions are the same as in Case I except that  $n_0$  observations, viz.,  $x_{01}, \dots, x_{0n_0}$  are taken on  $\gamma_0$ .

For selection procedure  $\gamma_6^{(2)}$ , the inequalities defining the procedure and corresponding to  $\bar{X}_{i;k} - c_{i;k} \gamma_0$  ( $i = 1, \dots, k$ ) are replaced by  $\bar{X}_{i;k} - c_{i;k}^{(2)} \gamma_0$  ( $i = 1, \dots, k$ ), respectively. The equation determining  $c_{i;k}^{(2)}$  is obtained as before and is given by

$$\int_0^\infty P(U_i - c_{i;k}^{(2)} \gamma_0) dt = 1, \quad (1.4.9)$$

where  $P(U_i < t)$  is the survival function in Theorem 1.4.1,  $\gamma_0$  the r.d.f. of  $\gamma_0$  and population  $\gamma_0$ .

If population  $\gamma_0$  has gamma distribution with density  $\phi(x; \gamma_0, \gamma_0)$  ( $\gamma_0$  known and  $\gamma_0$  unknown), then  $\phi(x; \gamma_0, \gamma_0; n_0, \gamma_0, \gamma_0)$ .

For selection procedure  $\gamma_7^{(2)}$ , the inequality defining the procedure is,

$$\beta_{k+1} < c^* \beta_0$$

and it can be shown  $c^* = \frac{\beta_0}{k\beta_0}$ .

For selection procedure  $\frac{\beta_0}{k}$ , the inequality defining the prior estimate and corresponding the  $(k-i)$ th step is

$$\frac{x_i - \beta_0}{x_i - \beta_0} < \frac{\beta_0}{\beta_0} \quad \text{where} \quad \beta_0 = \beta_0(u)$$

The equation determining  $c$  of  $\frac{\beta_0}{k}$  is given by

$$\int_0^{t_{k-i}} \int_0^{t_{k-i}} \frac{e^{-ut} e^{-ut}}{u^2} \frac{e^{-ut}}{u^2} du dt = \frac{e^{-2t}}{2} dt = p^*, \quad (1.1.1)$$

For selection procedure  $\frac{\beta_0}{k}$ , the inequality defining the prior estimate is

$$\frac{x_i - \beta_0}{x_i - \beta_0} < \frac{\beta_0}{\beta_0}, \quad (1.1.2)$$

and the equation determining  $c$  is given by Gupta and Soper (1970) as follows:

$$\int_0^{t_{k-i}} \int_0^{t_{k-i}} \frac{e^{-ut} e^{-ut}}{u^2} \frac{e^{-ut}}{u^2} du dt = \frac{e^{-2t}}{2} dt = p^*, \quad (1.1.3)$$

### 1.5. Selection Rules for the Location Parameter (prior estimate)

#### Ordering Prior Assumption

Since that we have only a partial information about the location parameter, that is, type of parameter space

$\mathbb{R}^k$  is a space for a set of  $k$  elements.

For a given  $k$ , the  $k$ -tuple  $(x_1, x_2, \dots, x_k)$  is a point in  $\mathbb{R}^k$ .

Let  $\{x_1, x_2, \dots, x_k\}$  be a set of  $k$  elements, and let  $\{B_1, B_2, \dots, B_k\}$  be  $k$  disjoint sets, say  $B_1, \dots, B_k$ , so that  $x_1 \in B_1, x_2 \in B_2, \dots, x_k \in B_k$ .

and for each  $b \in B_1, \dots, B_k$ , there is a unique  $x \in B_b$  such that  $b \in B_x$ .  
There is no order relation among the elements  $x_1, x_2, \dots, x_k$  in  $B_b$ .

Let  $b_i \in B_{x_i}$ , the number of elements in  $B_{x_i}$ , then we have

$$\frac{b_1}{b_2} < \frac{b_2}{b_3} < \dots < \frac{b_k}{b_1}$$

If we denote the new parameter  $p_i = b_i/b_{i+1}$ , then  $p_1, \dots, p_k$  are called the  $k$ -tuple parameter vector of  $x_1, \dots, x_k$ . We can also write  $p = (p_1, \dots, p_k)$  in an induced partial order.

Example. Suppose  $k = 3$ , and we have a partial order on  $\mathbb{R}^3$  given by  $x_1 < x_2 < x_3$ , and  $x_1 < x_3$ , and  $x_2 < x_3$ . We denote the elements  $x_1, x_2, x_3$  by  $a, b, c$  respectively. Then we can represent this partial order as in Figure 1.

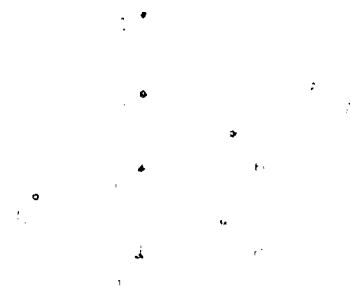


Figure 1. A partial order on  $\mathbb{R}^3$ .

then we have an induced partial order  $\preceq$  on  $\{1, 2, \dots, 10\}$  as in Figure 2.

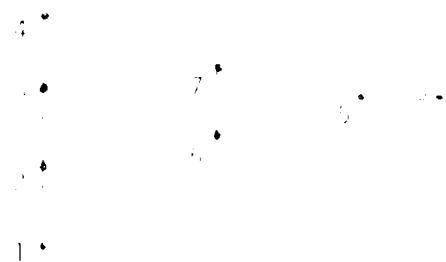


Figure 2. Induced partial order.

Now

$$B_0 = \{1, 2, 3, 4\}$$

$$B_1 = \{1, 2, 3, 4, 5, 6\}$$

$$B_2 = \{1, 2, 3, 4, 5, 6, 7\}$$

It is clear that the induced partial order is not unique. For example, we can partition  $\{1, 2, \dots, 10\}$  into three other subsets  $B_0^1, B_1^1, B_2^1$  where

$$B_0^1 = \{1, 2, 3, 4\}$$

$$B_1^1 = \{1, 2, 3, 4, 5, 6, 7\}$$

$$B_2^1 = \{8, 9, 10\}$$

From the local to global ordering, a selection procedure  $\{B_0, B_1, B_2\}$  is defined as follows:

Definition 1.1.1. We get the selection procedure  $\{B_0, B_1, B_2\}$  as follows:

Suppose  $\{B_0, B_1, B_2\}$  are the incident property and that the elements of  $B_0, B_1, B_2$  are the incident property of the incident property.

selection procedure — for each subset  $\mathcal{S}_j$  we have to find the maximum  $\hat{P}_{\mathcal{S}_j}^{\text{opt}}$

probability of a correct selection is not less than  $1 - \alpha_j^{\text{opt}}$

subset  $\mathcal{S}_j$  we may use selection procedure  $\hat{P}_{\mathcal{S}_j}^{\text{opt}} \geq \hat{P}_{\mathcal{S}_j}^{\text{opt}}$  with  $\hat{P}_{\mathcal{S}_j}^{\text{opt}}$

numerical Example. The selection procedure  $\hat{P}_{\mathcal{S}_j}^{\text{opt}}$  is given in the following

example

$$\begin{aligned} \inf_{\mathcal{S}_j} \hat{P}_{\mathcal{S}_j}^{\text{opt}} & (\text{CS}_j^{\text{opt}}) \\ & \text{such that} \\ & \inf_{\mathcal{S}_j \in \mathcal{S}_j^{\text{opt}}} \hat{P}_{\mathcal{S}_j}^{\text{opt}} \geq 1 - \alpha_j^{\text{opt}} \\ & \text{for } j = 1, 2, \dots, k \\ & \hat{P}_{\mathcal{S}_j}^{\text{opt}} = \hat{P}_{\mathcal{S}_j}^{\text{opt}}(b_j) \\ & \hat{P}_{\mathcal{S}_j}^{\text{opt}} = \hat{P}_{\mathcal{S}_j}^{\text{opt}}(b_j) \end{aligned}$$

where  $b_j$  is the parameter of a correlation with the subset  $\mathcal{S}_j$ .

**Remark 1.5.1.** For the selection problem in the case of one-dimensional distributions with respect to a certain criterion the following selection procedure can be used as a corollary from the previous theorem.

### 1.6. Comparison of the Bayesian procedure with others

#### 1.6.1. The location parameters of normal distribution

In this section we use Monte Carlo technique to compare the performance of the selection procedure (1.5.1) and (1.5.2) with the

$k$  independent populations, each population with distribution  $N(\mu_j, \sigma^2)$  with common known variance  $\sigma^2$  and common sample size  $n_j$ . Assume that the mean  $\mu_0$  of the control is known; without loss of generality we assume that  $\mu_0 = 0$  and  $\sigma/\sqrt{n} = 1$ .

In the simulation, we use Rubin and Hinkley's PVP-Random Variable Package, Purdue University Computing Center, to generate random numbers. For each  $k$ , we generated one random number (variable) for each population, then applied each selection procedure separately and repeated it ten thousand times; we used the relative frequency as an approximation of the exact values of the associated performance characteristics for each procedure. In Table V we use the 'old' wine notations:

$(\cdot, \dots, \cdot)$ ,  $\cdot_j$  is the parameter of population  $j$ .

$$PS = P(C_5)$$

PI = P(correctly rejecting all bad populations)

PC = P(correct classification of all population)

where the correct classification means that we select all the populations and reject all the impure populations.

(1) Expected number (size) of bad populations contained in the selected subset.

### 3.3 Expected size of the selected subset

Table V. Comparison of four procedures, namely, the  $\hat{\pi}_1$  procedure,  $\hat{\pi}_2$ ,  $\hat{\pi}_3$ ,  $\hat{\pi}_4$ , for each value of  $k$ , we assume that the true parameter  $\theta$  is the population is the one and only one parameter that with probability at least  $1 - \alpha$  is less than the control  $\theta_0 = 1$ . A glance at the table of  $\hat{\pi}_1$  in Table V shows that the performance can roughly be ordered as follows:

$$\hat{\pi}_1 > \hat{\pi}_3 > \hat{\pi}_2 > \hat{\pi}_4$$

i.e., procedure  $\hat{\pi}_1$  is the best one, followed by  $\hat{\pi}_3$ ,  $\hat{\pi}_2$  and  $\hat{\pi}_4$  are very close and both are better than  $\hat{\pi}_1$ . The ordering is based on the characteristics  $(\hat{\pi}_1, \hat{\pi}_2, \hat{\pi}_3, \hat{\pi}_4)$  of the four procedures given in Section 1 and 2. As the number of modulations  $k$  increases, the time to decide the true parameter and the three additional populations decreases. For the three parameters  $\lambda_1$ ,  $\lambda_2$  and  $\lambda_3$ , respectively, we find that  $\hat{\pi}_1$  is the best one, followed by  $\hat{\pi}_3$ ,  $\hat{\pi}_2$  and  $\hat{\pi}_4$  in  $\{1, 2, 3, 4\}$  in  $0.0164$ ,  $0.0144$ ,  $0.0134$  and  $0.0124$  respectively. This means that when  $k$  is even, and the additional populations are  $\lambda_1$ ,  $\lambda_2$  and  $\lambda_3$ , then  $\hat{\pi}_1$  is the best one, followed by  $\hat{\pi}_3$ ,  $\hat{\pi}_2$  and  $\hat{\pi}_4$ .

Table VI has the same structure as Table V, but for  $k = 1$ . In this case, we have two binary populations and one triple population. The performance characteristics of  $\hat{\pi}_1$ ,  $\hat{\pi}_2$ ,  $\hat{\pi}_3$  and  $\hat{\pi}_4$  are the same as in Table V, except that the time to decide the true parameter is longer.

In Table VII we compare the performance of the four procedures for  $k = 2$  and  $k = 3$  with the performance of the  $\hat{\pi}_1$  procedure for  $k = 1$ . The results are given in the last three rows of Table VII. The results show that the performance of the  $\hat{\pi}_1$  procedure is better than that of the other three procedures. The performance of the  $\hat{\pi}_1$  procedure is better than that of the other three procedures by about 30% for  $k = 1$  and for  $k = 2$  the performance of the  $\hat{\pi}_1$  procedure is better than that of the other three procedures by about 10%.

In Table 7.4 we compare the performance of the two models on the population of 1500 individuals. The results are very similar to those in Table 7.3, and we find that performance is not affected by the size of the population. The same result is obtained when we use the same model on the same results as before.

In Table V, we see that the only one which is not correctly (i.e., the true effect) is the one by  $\hat{F}_{\text{PE}}^{(1)}(x, y, t)$ . The effect of  $\hat{F}_{\text{PE}}^{(1)}(x, y, t)$  on PE, PPE and PDE have been estimated by  $\hat{F}_{\text{PE}}^{(1)}(x, y, t)$  and  $\hat{F}_{\text{PE}}^{(2)}(x, y, t)$ . The estimate of  $\hat{F}_{\text{PE}}^{(1)}(x, y, t)$  is the best. If we compare the two cases, we see that  $\hat{F}_{\text{PE}}^{(1)}(x, y, t)$  does not change the order of the terms in the equations.  $\hat{F}_{\text{PE}}^{(2)}(x, y, t)$  is the second best. For example, in a model  $y_i = \beta_0 + \beta_1 x_i + \beta_2 z_i + \varepsilon_i$ , the estimate of  $\beta_1$  is given by  $\hat{\beta}_1 = \hat{F}_{\text{PE}}^{(1)}(x, y, t) \hat{\beta}_1$  and the estimate of  $\beta_2$  is given by  $\hat{\beta}_2 = \hat{F}_{\text{PE}}^{(2)}(x, y, t) \hat{\beta}_2$ . In general,  $\hat{F}_{\text{PE}}^{(1)}(x, y, t)$  is better than  $\hat{F}_{\text{PE}}^{(2)}(x, y, t)$ .

Figure 1. Problem presentation for the 1000, 2000, and 3000 ms conditions. The 1000 ms condition is shown for reference.

The first question concerns the effect of the  $\alpha$  parameter on the  $\beta$  parameter. The second question concerns the effect of the  $\beta$  parameter on the  $\alpha$  parameter.

population has exponential distribution for each individual, and with unknown scale parameter  $\mu$  (see, e.g.,

In the simulation study, the parameter values used were those of Rubin and Stirkle's (1979) and the initial values were set to zero. The variables were specified as constant (1) and the results are based on two thousand sets of initial conditions. The results in Table VII and Table VIII are the same as in Table VI but the meaning of a bad population is that it is not a good population but a parameter which is greater than one (assumed one).

A good population means that its parameter is 100 times closer to one. The results of Table VI, I and Table VI, II and III show that we have the performance  $\rho_{\text{opt}}(\rho_{\text{opt}}) \approx \rho_{\text{opt}}(\rho_{\text{opt}})$  and that  $\rho_{\text{opt}}$  is slightly better than  $\rho_{\text{opt}}$ . There is only one fact that

Table of  $d_{13k}^{(1)}$  values, calculated by the direct and the indirect methods, to carry out the procedure  $\pi_1$  for the normalized multivariate  $\beta$  (the 1st order regression).

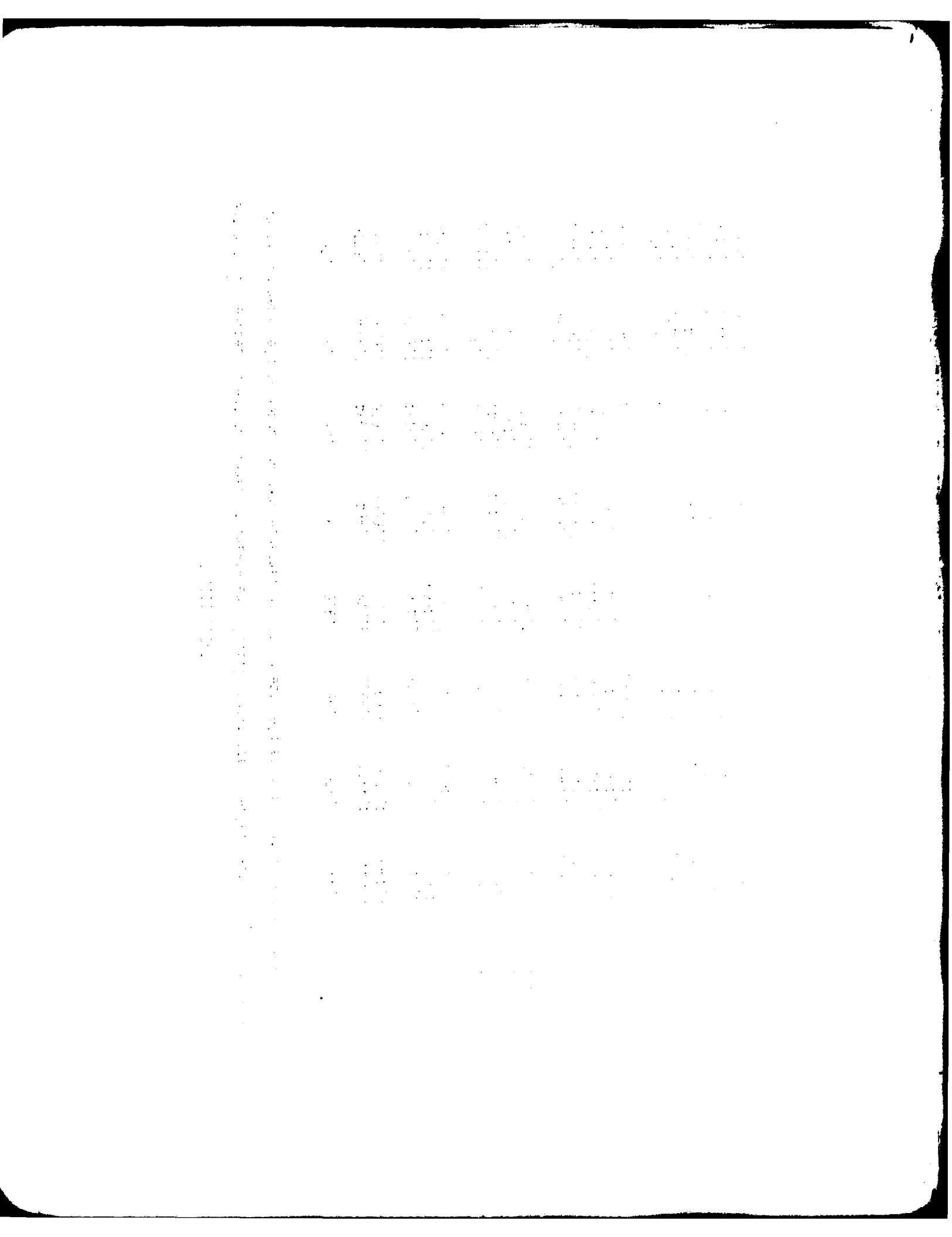
$k$	$d_{13k}^{(1)}$				
	$\alpha = 0.05$	$\alpha = 0.01$	$\alpha = 0.001$	$\alpha = 0.0001$	$\alpha = 0.00001$
1	2.13764	1.1669	0.74489	0.44346	0.27448
2	2.13347	1.1775	0.75720	0.45876	0.28447
3	2.13329	1.1787	0.76447	0.46966	0.28539
4	—	—	0.77152	0.47691	0.29570
5	—	—	0.77356	0.47963	0.29671
6	—	—	—	—	0.29772
7	1.3340	1.19787	0.78574	0.48496	0.30873
$d_{13k}^{(1)}$					
1	1.326	1.060	0.731	0.470	0.276
2	1.3264	1.0616	0.735	0.471	0.277
3	1.32736	1.0627	0.736	0.472	0.278
4	1.32735	1.0628	0.737	0.473	0.279
5	1.32736	1.0629	0.738	0.474	0.280
6	1.32736	1.0630	0.739	0.475	0.281
7	1.32736	1.0631	0.740	0.476	0.282
8	1.32736	1.0632	0.741	0.477	0.283
9	1.32736	1.0633	0.742	0.478	0.284
10	1.32736	1.0634	0.743	0.479	0.285

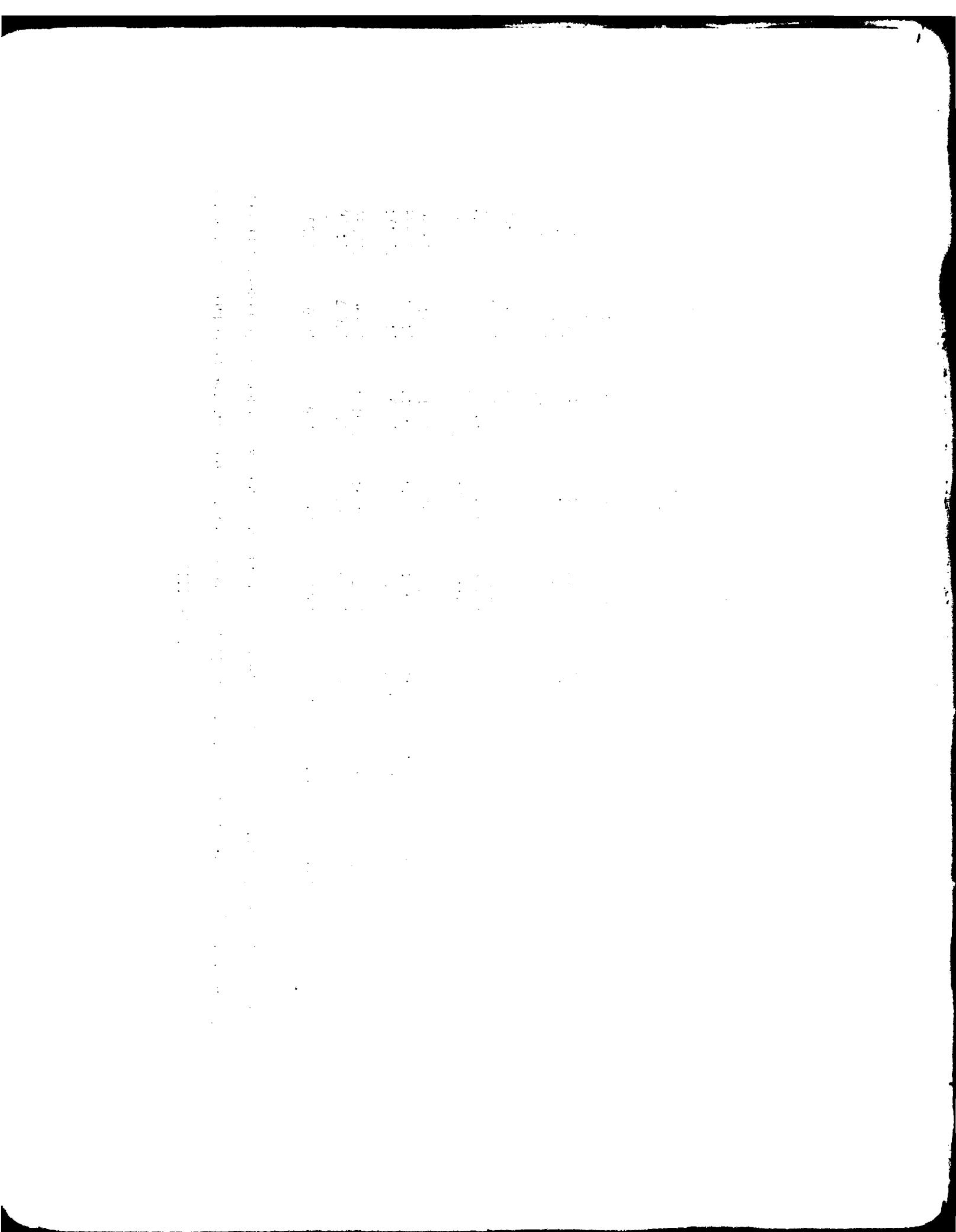
Table of  $d_{13k}^{(1)}$  values, calculated by the direct and the indirect methods, to carry out the procedure  $\pi_1$  for the normalized multivariate  $\beta$  (the 1st order regression).

1978, 1979)

Table of  $\alpha_{\text{per}}^{(2)}$  Additive correction to the  
procedure  $\alpha_{\text{per}}^{(1)}$  for the normal case of the model

$\alpha_{\text{per}}^{(2)}$	$\alpha_{\text{per}}^{(1)}$	$\alpha_{\text{per}}^{(2)}$	$\alpha_{\text{per}}^{(1)}$	$\alpha_{\text{per}}^{(2)}$	$\alpha_{\text{per}}^{(1)}$
0.00	0.00	0.00	0.00	0.00	0.00
0.05	0.05	0.05	0.05	0.05	0.05
0.10	0.10	0.10	0.10	0.10	0.10
0.15	0.15	0.15	0.15	0.15	0.15
0.20	0.20	0.20	0.20	0.20	0.20
0.25	0.25	0.25	0.25	0.25	0.25
0.30	0.30	0.30	0.30	0.30	0.30
0.35	0.35	0.35	0.35	0.35	0.35
0.40	0.40	0.40	0.40	0.40	0.40
0.45	0.45	0.45	0.45	0.45	0.45
0.50	0.50	0.50	0.50	0.50	0.50
0.55	0.55	0.55	0.55	0.55	0.55
0.60	0.60	0.60	0.60	0.60	0.60
0.65	0.65	0.65	0.65	0.65	0.65
0.70	0.70	0.70	0.70	0.70	0.70
0.75	0.75	0.75	0.75	0.75	0.75
0.80	0.80	0.80	0.80	0.80	0.80
0.85	0.85	0.85	0.85	0.85	0.85
0.90	0.90	0.90	0.90	0.90	0.90
0.95	0.95	0.95	0.95	0.95	0.95
1.00	1.00	1.00	1.00	1.00	1.00









the value  $\rho_{\text{max}} = 0.8$  in the first column of Table 1, we can see that the error rate  $\rho_{\text{err}}$  for the quantile-based procedure is within the expected range for the confidence interval.

Table 1

n	rho_max			
	0.1	0.2	0.3	0.4
10	0.0513	0.1000	0.1500	0.2000
20	0.0213	0.0400	0.0600	0.0800
30	0.0133	0.0200	0.0300	0.0400
40	0.0083	0.0100	0.0100	0.0100
50	0.0050	0.0050	0.0050	0.0050
60	0.0033	0.0033	0.0033	0.0033
70	0.0021	0.0021	0.0021	0.0021
80	0.0013	0.0013	0.0013	0.0013
90	0.00083	0.00083	0.00083	0.00083
100	0.00050	0.00050	0.00050	0.00050
150	0.00021	0.00021	0.00021	0.00021
200	0.00013	0.00013	0.00013	0.00013
300	0.000083	0.000083	0.000083	0.000083
400	0.000050	0.000050	0.000050	0.000050
500	0.000033	0.000033	0.000033	0.000033
600	0.000021	0.000021	0.000021	0.000021
700	0.000013	0.000013	0.000013	0.000013
800	0.0000083	0.0000083	0.0000083	0.0000083
900	0.0000050	0.0000050	0.0000050	0.0000050
1000	0.0000033	0.0000033	0.0000033	0.0000033

Table 1. The quantile-based procedure for the confidence interval of the mean.

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1		2		3		4		5		6	
1	1	2	2	3	3	4	4	5	5	6	6
2	2	3	3	4	4	5	5	6	6	7	7
3	3	4	4	5	5	6	6	7	7	8	8
4	4	5	5	6	6	7	7	8	8	9	9
5	5	6	6	7	7	8	8	9	9	10	10
6	6	7	7	8	8	9	9	10	10	11	11
7	7	8	8	9	9	10	10	11	11	12	12
8	8	9	9	10	10	11	11	12	12	13	13
9	9	10	10	11	11	12	12	13	13	14	14
10	10	11	11	12	12	13	13	14	14	15	15
11	11	12	12	13	13	14	14	15	15	16	16
12	12	13	13	14	14	15	15	16	16	17	17
13	13	14	14	15	15	16	16	17	17	18	18
14	14	15	15	16	16	17	17	18	18	19	19
15	15	16	16	17	17	18	18	19	19	20	20
16	16	17	17	18	18	19	19	20	20	21	21
17	17	18	18	19	19	20	20	21	21	22	22
18	18	19	19	20	20	21	21	22	22	23	23
19	19	20	20	21	21	22	22	23	23	24	24
20	20	21	21	22	22	23	23	24	24	25	25
21	21	22	22	23	23	24	24	25	25	26	26
22	22	23	23	24	24	25	25	26	26	27	27
23	23	24	24	25	25	26	26	27	27	28	28
24	24	25	25	26	26	27	27	28	28	29	29
25	25	26	26	27	27	28	28	29	29	30	30
26	26	27	27	28	28	29	29	30	30	31	31
27	27	28	28	29	29	30	30	31	31	32	32
28	28	29	29	30	30	31	31	32	32	33	33
29	29	30	30	31	31	32	32	33	33	34	34
30	30	31	31	32	32	33	33	34	34	35	35
31	31	32	32	33	33	34	34	35	35	36	36
32	32	33	33	34	34	35	35	36	36	37	37
33	33	34	34	35	35	36	36	37	37	38	38
34	34	35	35	36	36	37	37	38	38	39	39
35	35	36	36	37	37	38	38	39	39	40	40
36	36	37	37	38	38	39	39	40	40	41	41
37	37	38	38	39	39	40	40	41	41	42	42
38	38	39	39	40	40	41	41	42	42	43	43
39	39	40	40	41	41	42	42	43	43	44	44
40	40	41	41	42	42	43	43	44	44	45	45
41	41	42	42	43	43	44	44	45	45	46	46
42	42	43	43	44	44	45	45	46	46	47	47
43	43	44	44	45	45	46	46	47	47	48	48
44	44	45	45	46	46	47	47	48	48	49	49
45	45	46	46	47	47	48	48	49	49	50	50
46	46	47	47	48	48	49	49	50	50	51	51
47	47	48	48	49	49	50	50	51	51	52	52
48	48	49	49	50	50	51	51	52	52	53	53
49	49	50	50	51	51	52	52	53	53	54	54
50	50	51	51	52	52	53	53	54	54	55	55
51	51	52	52	53	53	54	54	55	55	56	56
52	52	53	53	54	54	55	55	56	56	57	57
53	53	54	54	55	55	56	56	57	57	58	58
54	54	55	55	56	56	57	57	58	58	59	59
55	55	56	56	57	57	58	58	59	59	60	60
56	56	57	57	58	58	59	59	60	60	61	61
57	57	58	58	59	59	60	60	61	61	62	62
58	58	59	59	60	60	61	61	62	62	63	63
59	59	60	60	61	61	62	62	63	63	64	64
60	60	61	61	62	62	63	63	64	64	65	65
61	61	62	62	63	63	64	64	65	65	66	66
62	62	63	63	64	64	65	65	66	66	67	67
63	63	64	64	65	65	66	66	67	67	68	68
64	64	65	65	66	66	67	67	68	68	69	69
65	65	66	66	67	67	68	68	69	69	70	70
66	66	67	67	68	68	69	69	70	70	71	71
67	67	68	68	69	69	70	70	71	71	72	72
68	68	69	69	70	70	71	71	72	72	73	73
69	69	70	70	71	71	72	72	73	73	74	74
70	70	71	71	72	72	73	73	74	74	75	75
71	71	72	72	73	73	74	74	75	75	76	76
72	72	73	73	74	74	75	75	76	76	77	77
73	73	74	74	75	75	76	76	77	77	78	78
74	74	75	75	76	76	77	77	78	78	79	79
75	75	76	76	77	77	78	78	79	79	80	80
76	76	77	77	78	78	79	79	80	80	81	81
77	77	78	78	79	79	80	80	81	81	82	82
78	78	79	79	80	80	81	81	82	82	83	83
79	79	80	80	81	81	82	82	83	83	84	84
80	80	81	81	82	82	83	83	84	84	85	85
81	81	82	82	83	83	84	84	85	85	86	86
82	82	83	83	84	84	85	85	86	86	87	87
83	83	84	84	85	85	86	86	87	87	88	88
84	84	85	85	86	86	87	87	88	88	89	89
85	85	86	86	87	87	88	88	89	89	90	90
86	86	87	87	88	88	89	89	90	90	91	91
87	87	88	88	89	89	90	90	91	91	92	92
88	88	89	89	90	90	91	91	92	92	93	93
89	89	90	90	91	91	92	92	93	93	94	94
90	90	91	91	92	92	93	93	94	94	95	95
91	91	92	92	93	93	94	94	95	95	96	96
92	92	93	93	94	94	95	95	96	96	97	97
93	93	94	94	95	95	96	96	97	97	98	98
94	94	95	95	96	96	97	97	98	98	99	99
95	95	96	96	97	97	98	98	99	99	100	100

a<sub>1</sub> (1)

1	1.6702	1.5958	1.5713	1.4976	1.4410	1.3717	
2	1.7039	1.6668	1.6367	1.4596	1.3980	1.3217	
3	1.7406	1.6564	1.6147	1.4561	1.3728	1.3007	
4	1.7440	1.6976	1.6537	1.4579	1.3711	1.3044	
5	1.7451	1.6895	1.6556	1.4584	1.3704	1.3031	
6	1.7455	1.6888	1.6557	1.4585	1.3707	1.3035	
7	1.7456	1.6879	1.6557	1.4586	--	--	
8	1.7457	--	--	--	--	--	
10	--	--	--	--	--	--	
1	2.79437	2.7567	2.7615	2.7397	2.7479	2.7444	
2	3.0126	2.8194	2.7951	2.6779	2.7136	2.7433	
3	3.0264	2.8223	2.7143	2.6396	2.7144	2.7434	
4	3.0301	2.8320	2.7156	2.6368	2.7156	2.7443	
5	3.0313	2.8328	2.7162	2.6374	2.7162	2.7446	
6	3.0317	2.8331	2.7164	2.6376	2.7164	2.7447	
7	3.0318	2.8337	2.7166	2.6377	--	--	
8	3.0319	--	--	--	--	--	
10	--	--	--	--	--	--	
1	4.1496	4.1134	4.1164	3.7741	3.6258	3.5617	
2	4.1736	4.1168	4.1167	3.7686	3.6196	3.5549	
3	4.1744	4.1096	4.1086	3.7717	3.6269	3.5636	
4	4.17470	4.1099	4.1086	3.7737	3.6274	3.5641	
5	4.17491	4.1101	4.1166	3.7744	3.6277	3.5643	
6	4.17496	4.1114	4.1167	3.7746	3.6278	3.5645	
7	4.17498	4.1114	4.1167	3.7747	--	--	
10	--	--	--	--	--	--	

1. *Leucosia* *leucosia* (L.) *leucosia* (L.)

2. *Leucosia* *leucosia* (L.) *leucosia* (L.)

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13. *Leucosia* *leucosia* (L.) *leucosia* (L.)

14. *Leucosia* *leucosia* (L.) *leucosia* (L.)

15. *Leucosia* *leucosia* (L.) *leucosia* (L.)

16. *Leucosia* *leucosia* (L.) *leucosia* (L.)

17. *Leucosia* *leucosia* (L.) *leucosia* (L.)

18. *Leucosia* *leucosia* (L.) *leucosia* (L.)

19. *Leucosia* *leucosia* (L.) *leucosia* (L.)

20. *Leucosia* *leucosia* (L.) *leucosia* (L.)

should be used to indicate the presence of a significant difference between the two groups, provided that the two groups are not too dissimilar in size.

$k = 3, \quad (-1, 1, 1)$	$\{ \}$	$\{ \}$	$\{ \}$	$\{ \}$
P1	1.9453	1.0000	1.0000	1.0000
P2	1.3854	1.0000	1.0000	1.0000
P3	1.3507	1.0000	1.0000	1.0000
P4	1.6146	1.0000	1.0000	1.0000
P5	1.6146	1.0000	1.0000	1.0000
P6	1.5599	1.0000	1.0000	1.0000
$k = 4, \quad (-1, 0, 1)$	$\{ \}$	$\{ \}$	$\{ \}$	$\{ \}$
P1	1.6531	1.0000	1.0000	1.0000
P2	1.5744	1.0000	1.0000	1.0000
P3	1.5744	1.0000	1.0000	1.0000
P4	1.6151	1.0000	1.0000	1.0000
P5	1.6151	1.0000	1.0000	1.0000
P6	1.5774	1.0000	1.0000	1.0000
$k = 4, \quad (-1, 0, 1, -1)$	$\{ \}$	$\{ \}$	$\{ \}$	$\{ \}$
P1	1.9580	1.0000	1.0000	1.0000
P2	1.9564	1.0000	1.0000	1.0000
P3	1.8744	1.0000	1.0000	1.0000
P4	1.6136	1.0000	1.0000	1.0000
P5	1.6136	1.0000	1.0000	1.0000
P6	1.5590	1.0000	1.0000	1.0000
$k = 5, \quad (-1, 0, 1, -1, 1)$	$\{ \}$	$\{ \}$	$\{ \}$	$\{ \}$
P1	1.6534	1.0000	1.0000	1.0000
P2	1.5745	1.0000	1.0000	1.0000
P3	1.5745	1.0000	1.0000	1.0000
P4	1.6152	1.0000	1.0000	1.0000
P5	1.6152	1.0000	1.0000	1.0000
P6	1.5775	1.0000	1.0000	1.0000

Simulation results show that the proposed method can obtain the accurate estimation of the initial state and input, and the proposed method is more robust than the LQR method.

For more information, contact the U.S. Environmental Protection Agency's Office of Water, Washington, D.C. 20460, (202) 265-2600.

1. *What is the name of the author of the book?*  
2. *What is the name of the book?*  
3. *What is the name of the publisher?*

4. *What is the name of the author of the book?*  
5. *What is the name of the book?*  
6. *What is the name of the publisher?*

7. *What is the name of the author of the book?*  
8. *What is the name of the book?*  
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10. *What is the name of the author of the book?*  
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12. *What is the name of the publisher?*

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14. *What is the name of the book?*  
15. *What is the name of the publisher?*

16. *What is the name of the author of the book?*  
17. *What is the name of the book?*  
18. *What is the name of the publisher?*

19. *What is the name of the author of the book?*  
20. *What is the name of the book?*  
21. *What is the name of the publisher?*

22. *What is the name of the author of the book?*  
23. *What is the name of the book?*  
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25. *What is the name of the author of the book?*  
26. *What is the name of the book?*  
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28. *What is the name of the author of the book?*  
29. *What is the name of the book?*  
30. *What is the name of the publisher?*

31. *What is the name of the author of the book?*  
32. *What is the name of the book?*  
33. *What is the name of the publisher?*

34. *What is the name of the author of the book?*  
35. *What is the name of the book?*  
36. *What is the name of the publisher?*

37. *What is the name of the author of the book?*  
38. *What is the name of the book?*  
39. *What is the name of the publisher?*

40. *What is the name of the author of the book?*  
41. *What is the name of the book?*  
42. *What is the name of the publisher?*

43. *What is the name of the author of the book?*  
44. *What is the name of the book?*  
45. *What is the name of the publisher?*

TABLE 4

Simulation results for the comparative performance of various statistical procedures for the model mean  $\mu = (\mu_1, \mu_2, \mu_3)$  (not including the constraint  $\mu_1 + \mu_2 + \mu_3 = 1$ ) under simple symmetric priors.

$\mu = (0.3, 0.3, 0.4)$			
$\hat{\mu}_1$	0.3000	0.3000	0.3000
PS	1.0000	0.9993	0.9993
PI	1.0013	0.9999	0.9999
PI	1.0013	0.9999	0.9999
PI	1.0007	0.9999	0.9999
PS	1.0001	0.9999	0.9999
PS	1.0007	0.9999	0.9999
$\mu = (0.3, 0.4, 0.3)$			
$\hat{\mu}_1$	0.3000	0.3000	0.3000
PS	1.0000	0.9993	0.9993
PI	1.0017	0.9999	0.9999
PI	1.0021	0.9999	0.9999
PI	1.0010	0.9999	0.9999
PS	1.0007	0.9999	0.9999
PS	1.0007	0.9999	0.9999
$\mu = (0.4, 0.3, 0.3)$			
$\hat{\mu}_1$	0.3000	0.3000	0.3000
PS	1.0000	0.9993	0.9993
PI	1.0012	0.9999	0.9999
PI	1.0019	0.9999	0.9999
PI	1.0013	0.9999	0.9999
PS	1.0007	0.9999	0.9999
PS	1.0007	0.9999	0.9999

Table V, 3

Simulation results for the comparative performance of various selection procedures for the normal means problem (notations explained in Section 1.6.1) under simple ordering prior.

$\mu^* = 900$

	$k = 2, \gamma = (-1, +1)$	$k = 3, \gamma = (-1, +1, +1)$	$k = 4, \gamma = (-1, +1, +1, +1)$	$k = 5, \gamma = (-1, +1, +1, +1, +1)$
P5	1.0000	1.0000	1.0000	1.0000
PI	1.5405	1.5343	1.5337	1.5272
PC	1.5405	1.5343	1.5337	1.5272
U1	1.8331	1.8337	1.8351	1.8394
U3	1.7116	1.7349	1.7432	1.7573
US	1.8331	1.8337	1.8351	1.8394
P5	1.9957	1.9957	1.9957	1.9957
PI	1.5365	1.5343	1.5337	1.5272
PC	1.5397	1.5343	1.5337	1.5272
U1	1.8347	1.8363	1.8376	1.8394
U3	1.7357	1.7341	1.7415	1.7491
US	1.8279	1.8330	1.8373	1.8394
P5	1.9924	1.9924	1.9924	1.9924
PI	1.5274	1.5263	1.5264	1.5267
PC	1.5197	1.5263	1.5267	1.5272
U1	1.8403	1.8390	1.8385	1.8377
U3	1.7055	1.7072	1.7142	1.7276
US	1.8390	1.8370	1.8372	1.8374
P5	1.9901	1.9901	1.9901	1.9901
PI	1.5217	1.5207	1.5207	1.5207
PC	1.5177	1.5207	1.5207	1.5207
U1	1.8461	1.8461	1.8461	1.8461
U3	1.7151	1.7174	1.7216	1.7276
US	1.8341	1.8330	1.8373	1.8394

TABLE VII

Simulation results for the comparative performance of various statistical procedures for the gamma means problem (notation explained in Section 1.6.1) under simple order restriction.

$P^* = 0.90$

$\kappa = 4, \mu = \{.3, .6, .9, 1.2\}$

	$\mu = .3$	$\mu = .6$	$\mu = .9$	$\mu = 1.2$
PS	.0840	.0848	.0855	.0862
PI	.7629	.7745	.7872	.7997
PC	.1860	.1597	.1375	.1167
SI	.7980	.7955	.7935	.7917
ED	.0319	.0320	.0321	.0322
EF	2.7329	3.8190	4.0417	4.3063

$\kappa = 6, \mu = \{.3, .6, .9, 1.2, 1.5\}$

	$\mu = .3$	$\mu = .6$	$\mu = .9$	$\mu = 1.2$	$\mu = 1.5$
PS	.0905	.0910	.0915	.0920	.0925
PI	.6990	.6950	.6845	.6743	.6643
PC	.0695	.0750	.0815	.0880	.0945
SI	1.5040	1.4560	1.3870	1.3180	1.2500
ED	.1868	.1969	.2068	.2167	.2266
EF	4.4945	4.5470	4.5895	4.6320	4.6745

$\kappa = 4, \mu = \{.3, .6, .9, 1.2, 1.5\}$

	$\mu = .3$	$\mu = .6$	$\mu = .9$	$\mu = 1.2$	$\mu = 1.5$
PS	.0625	.0630	.0635	.0640	.0645
PI	.7625	.7667	.7710	.7753	.7795
PC	.2370	.2373	.2375	.2377	.2379
SI	1.0790	1.1110	1.1430	1.1750	1.2070
ED	.5304	.5310	.5317	.5324	.5331
EF	2.2475	2.3170	2.3865	2.4560	2.5255

$\kappa = 6, \mu = \{.3, .6, .9, 1.2, 1.5\}$

	$\mu = .3$	$\mu = .6$	$\mu = .9$	$\mu = 1.2$	$\mu = 1.5$
PS	.0600	.0600	.0600	.0600	.0600
PI	.7600	.7600	.7600	.7600	.7600
PC	.2310	.2310	.2310	.2310	.2310
SI	1.0610	1.1030	1.1350	1.1670	1.2000
ED	.5250	.5250	.5250	.5250	.5250
EF	2.2150	2.2850	2.3550	2.4250	2.4950

TABLE VI.

Simulation results for the comparative performance of various selection procedures for the gamma mean problem (notation explained in Section 1.6.1) under simple ordering prior.

P = .900							
k = 4, n = (1, 1, .6, 1, 1, 1, 1)		6		7		8	
PS		.9995		.9995		1.0000	
PI		.0690		.0645		.0545	
PC		.0685		.0640		.0545	
EI		1.4940		1.5345		1.5315	
EJ		.2120		.2750		.2204	
ES		3.4935		3.5340		3.5315	
k = 5, n = (1, 1, .6, 1, 1, 1, 1, 2)		6		7		8	
PS		.9990		.9995		1.0000	
PI		.0540		.0530		.0295	
PC		.0530		.0520		.0295	
EI		1.9205		1.9700		2.1385	
EJ		.6140		.6646		.7727	
ES		3.9195		3.9630		4.1385	
k = 4, n = (1, 1, 1, 4, 1, 1, 1, 1)		6		7		8	
PS		1.0000		1.0000		1.0000	
PI		.0375		.0375		.0375	
PC		.0375		.0375		.0375	
EI		.6890		.7475		.8115	
EJ		.6796		.7371		.8041	
ES		.6890		.7475		.8115	
k = 5, n = (1, 1, 1, 4, 1, 1, 1, 1, 2)		6		7		8	
PS		1.0000		1.0000		1.0000	
PI		.0315		.0315		.0315	
PC		.0315		.0315		.0315	
EI		.9730		.9730		.9600	
EJ		1.0000		1.0000		1.0000	
ES		1.0000		1.0000		1.0000	

CHAPTER II  
SAYE'S  $\star$  SELECTION RULES  
FOR SELECTING A SUBSET CONTAINING  
THE BEST POPULATION

2.1. Introduction

Suppose we have  $k$  ( $k \geq 2$ ) independent populations  $\pi_1, \pi_2, \dots, \pi_k$  and that the random variable  $x_{ij}$  associated with  $\pi_j$  has a distribution with  $\alpha$  known parameters  $\alpha_{ij}$ ,  $i = 1, 2, \dots, n_j$ . First, we give some definitions.

**Definition 2.1.1.** The population  $\pi_j$  is the best population of  $\pi_1, \pi_2, \dots, \pi_k$  for all  $j \neq i$ , if there are more than one population with this condition we arbitrarily tag one of them and call it the best population which is not the best, we called a non-best population.

Assume that we have  $n_j$  independent observations  $x_{ij}$ ,  $i = 1, 2, \dots, n_j$ , for population  $\pi_j$ ,  $j = 1, 2, \dots, k$ . Let  $\hat{\alpha}_{ij}$  be the estimate of  $\alpha_{ij}$ ,  $i = 1, 2, \dots, n_j$ , so that  $\hat{\alpha}_{ij}$  are independent. Let  $\hat{\alpha}_{ij}^*$  be the estimate of  $\alpha_{ij}$  for the best population  $\pi_j$ , for each  $i$ , we have  $\hat{\alpha}_{ij}^* = \hat{\alpha}_{ij}$  for  $i = 1, 2, \dots, n_j$ . Let  $\hat{\alpha}_{ij}^*$  be the estimate of  $\alpha_{ij}$  for the non-best population  $\pi_j$ , for  $i = 1, 2, \dots, n_j$ . Then we have

$$\hat{\alpha}_{ij}^* = \hat{\alpha}_{ij} \quad \text{for } i = 1, 2, \dots, n_j \text{ and } j = 1, 2, \dots, k.$$

The estimate  $\hat{\alpha}_{ij}^*$  is called the  $\star$  estimate of  $\alpha_{ij}$  for the non-best population  $\pi_j$ .

$\{1, 2, \dots, k\}$ . An action  $A \in \mathcal{A}$  is the selection of some subset of the  $k$  populations, i.e.  $i \in A$  means that  $i$  is included in the selected subset. An action  $A \in \mathcal{A}$  is called a correct selection (CS) if the best population is included in the selected subset  $A$ .

**Definition 2.1.2.** A measurable function  $\pi$  defined on  $\mathcal{X} \times \mathcal{A}$  is called a selection procedure provided that for each  $x \in \mathcal{X}$ , we have

$$0 \leq \pi(x, A)$$

and

$$\sum_{A \in \mathcal{A}} \pi(x, A) = 1$$

where  $\pi(x, A)$  denotes the probability that the subset  $A$  is selected when  $x$  is observed.

The individual selection probability  $\pi_i(x)$  for the population  $i$  is then given by

$$\pi_i(x) = \sum_{A \ni i} \pi(x, A)$$

where the summation is over all subsets  $A$  which contain  $i$ . The selection probability  $\pi_i(x)$  takes on only values 0 or 1, i.e. we say that the selection procedure  $\pi(x, A)$  is completely efficient if this is the case. It is called a non-randomized procedure.

**Definition 2.1.3.** Two selection procedures  $\pi$  and  $\pi'$  are equivalent if they have the same individual selection probabilities, i.e. if

$$\text{for all } x, i = 1, \dots, k,$$

Hence we can use the following definition, replacing  $\pi$  by  $\pi'$ :

**Definition 3.1.4.** A subset selection rule  $\tau$  is a measurable function from  $\mathbb{R}^k$  to  $\mathbb{R}^k$ ,

$$\tau(x) = (\tau_1(x), \dots, \tau_k(x))$$

with

$$0 \leq \tau_j(x) \leq 1, \quad j = 1, \dots, k.$$

If  $\tau_j$ 's are 0 or 1, the rule is nonstochastic. Note that, by definition 3.1.1, we have  $\sum_{j=1}^k \tau_j(x) \leq 1$ .

Suppose our goal is to find a nontrivial subset which contains the best population. A large body of literature exists in the area of subset selection procedures (see Gupta and Sanchezaram (1971), Gupta (1956, 1966) gave maximum-type subset selection procedures. Gupta (1972) studied the performance of Gupta-type maximum procedure, the real-type average procedure (Seal (1956, 1967)) and the Bayesian procedure. Berger (1971a) and Berger and Gupta (1970) proved that Gupta-type maximum procedure is minimax under certain loss functions. In the context of the set-in approach to the subset selection problem, Gold and Puri (1970), Chernoff and Sobay (1971), Buckley and Gray (1972), Gupta and Puri (1972) and Puri (1979) gave different formulation under different loss functions. The loss functions proposed by the last three consist of two parts: the linear and the quadratic loss parts. According to the nature of the loss function, one of the two is always dominant, whenever the other is not, one may wish to try some other method of attack.

On the other hand, an important family of subset selection procedures is the family of nonstochastic subset selection procedures. The most well-known member of this family is the  $\tau$  rule, which is defined as

the different sampling procedures can be derived in the same way. However, the exact sampling distributions of the different estimators usually have to wait until different data are obtained. In this chapter we focus on the properties of the estimators for the type models of homogeneity, and, although the effect of the exact sampling properties of the estimators is not explicitly discussed, we may wish to relate (modified) the so-called Bayesian prior.

In this chapter we define the posterior  $\hat{\pi}^B$  estimator in section 2.1. Two Bayesian selection procedure,  $\hat{\pi}_k^B$  and  $\hat{\pi}_s^B$  are proposed in section 2.2 and section 2.3 separately, and their properties are discussed in section 2.4. In section 2.5 we discuss their applications to gamma distributions. In section 2.6, procedure  $\hat{\pi}_s^B$  is compared with a classical selection procedure. An application for the problem of selecting the best  $k$   $\times$   $k$  block of the initial distributions  $\mathbf{M}_k$  is given in section 2.7. In section 2.8 we discuss their applications to the selection problems for Poisson distributions and beta distributions, and their relation to the selection of gamma distributions. In section 2.9, we deal with the comparison of the performance of selection procedures  $\hat{\pi}_k^B$  and  $\hat{\pi}_s^B$ , where  $\hat{\pi}_k^B$  and  $\hat{\pi}_s^B$  are the Bayesian selection procedures based on single prior and sample posterior, respectively (see Gupta and Bhattacharya (1986), Bhattacharya and Gupta (1987) and Bhattacharya and Gupta (1989)). The comparison is based on Monte Carlo simulation. In addition to the comparison in terms of the expected error and the variance of the estimator, the performance of the two procedures is evaluated in terms of the expected value of the sample posterior probability of the true distribution. The two procedures are also compared in terms of the expected value of the sample posterior probability of the true distribution.

## 2.2. Definitions of the Posterior- $P^*$ Condition and the Non-randomized Bayes- $P^*$ Procedure $\pi_{\text{B}}^*$

Let  $\{1, 2, \dots, k\}$  be the ordered unknown  $\gamma_1^{\text{true}}$  under consideration. If the prior distribution  $\pi$  for  $\gamma = (\gamma_1, \dots, \gamma_k)$  then the posterior probability of a correct selection under selection procedure  $\pi$ , given  $y = x$ , is

$$P(\text{CS}_{\pi}, y = x) = \sum_{i=1}^k p_i(x) \pi_i$$

where

$$p_i(x) = P(\gamma_i \text{ is the best} | y = x).$$

It is clear

$$\sum_{i=1}^k p_i(x) \leq 1.$$

**Definition 2.2.1.** Given a number  $P^* (0 < P^* \leq 1)$  and the prior  $\pi$ , we say a selection procedure  $\pi$  satisfies the posterior- $P^*$  condition if

$$P(\text{CS}_{\pi}, y = x) \geq P^* \quad \text{for all } x.$$

**Remark 2.2.1.** The posterior- $P^*$  condition is based on the prior distribution  $\pi$  and is different from the usual so-called  $P^*$ -condition.

**Definition 2.2.2.** The loss function  $\psi_{\pi}$  is defined by  $\psi_{\pi}(\gamma, y)$  where  $\gamma$  is the size (number) of populations associated with the best of  $\gamma$ , the loss function  $\psi_{\pi}$  is defined by  $\psi_{\pi}(\gamma, A) = \sum_{i \in A} \psi_{\pi}(\gamma_i, y)$

which is the number of the non-best populations selected by  $\pi$  (note that  $A \subseteq \{1, 2, \dots, k\}$ ).

Note that the indicator function

$$\mathbb{I}_{\{\gamma_i \in \gamma\}} = \begin{cases} 1 & \text{if } \gamma_i \in \gamma \\ 0 & \text{otherwise} \end{cases}$$

definition 2.2.3. Given a number  $P^* \frac{1}{k} \leq P^* \leq 1$  and the prior  $\pi$ , we define the class  $\omega_{NR}(\cdot, P^*)$  as follows.

$\omega_{NR}(\cdot, P^*) = \{\cdot\} \text{ is any non-randomized rule which satisfies the posterior-} P^* \text{ condition.}$

For the sake of convenience sometimes we will use  $\omega_{NP}$  instead of  $\omega_{NR}(\cdot, P^*)$ .

definition 2.2.4. Given a number  $P^* \frac{1}{k} \leq P^* \leq 1$ , a prior  $\pi$  and the loss function  $L$ , a selection procedure  $\omega_{NP}(\cdot, P^*)$  is called a non-randomized Bayes- $P^*$  procedure (rule) if  $\cdot$  is a Bayes rule in the class  $\omega_{NP}(\cdot, P^*)$ .

Let  $p_{(1)}(x) \leq \dots \leq p_{(k)}(x)$  be the ordered  $p_i(x)$ 's and  $\pi_{(i)}$  be the population associated with  $p_{(i)}(x)$ ,  $i = 1, \dots, k$ , then a subset selection rule  $\cdot$  is completely specified by  $\omega_{(1)}, \dots, \omega_{(k)}$ , where  $\omega_{(i)}$  is defined by

$$\omega_{(i)}(x) = P\{x \in \omega_i\}, \quad i = 1, \dots, k.$$

Next, we propose a non-randomized selection rule which belongs to

$\omega_{NP}(\cdot, P^*)$ .

definition 2.2.5. Given a number  $P^* \frac{1}{k} \leq P^* \leq 1$ , a prior  $\pi$  and a prior distribution  $\pi$ , the selection rule  $\omega_{NP}$  is defined by  $\omega_{NP}(x) = \omega(x)$ , where

$$\omega(x) = \begin{cases} 1, & \text{if } x \in \omega(x) \\ 0, & \text{otherwise} \end{cases}$$

and  $j(x)$  is the maximum integer such that

$$\sum_{i=1}^k p_{\{i\}}(x) < p^*.$$

Lemma 2.2.1.  $\epsilon_{NR}^B \in \mathcal{L}_{NR}^B$ .

Proof. Follows from the definition of  $\epsilon_{NR}^B$ .

Theorem 2.2.1. Given a number  $p_{\{k\}}^{*+1} < p^* < 1$ , the prior  $\epsilon$  and the function  $L_1$ , the selection procedure  $\epsilon_{NR}^B$  is a non-randomized Bayes rule.

Proof. It is sufficient to show that the selection procedure  $\epsilon_{NR}^B$  has the smallest posterior risk in the class  $\mathcal{L}_{NR}^B(\epsilon, p^*)$ . Given the observation  $X = x$ . Let the posterior risk of  $\epsilon \in \mathcal{L}_{NR}^B(\epsilon, p^*)$  be  $r(x, \epsilon)$ .

$$r(x, \epsilon_{NR}^B) = k - j + 1$$

and

$$\sum_{i=j+1}^k p_{\{i\}}(x) < p^*$$

for some  $j$ ,  $1 \leq j \leq k$ .

Hence the inequality

$$r(x, \epsilon) \geq r(x, \epsilon_{NR}^B)$$

is not true for any  $\epsilon \in \mathcal{L}_{NR}^B(\epsilon, p^*)$ ,  $\epsilon \neq \epsilon_{NR}^B$ . Therefore, the result follows.

Theorem 2.2.2. Theorem 2.2.1 also holds when we replace the loss  $L_1$  by

Proof. Under the loss function  $L_2$ , the posterior risk of the selection procedure  $\pi \in \mathcal{L}_{NR}^*(\pi, p^*)$  is

$$r(x, \pi) = \sum_{i=1}^k \pi(i)(x)[i - p_{[i]}(x)], \text{ given } x \in \mathcal{X}.$$

By Theorem 2.2.1, we have

$$\sum_{i=1}^k \pi_{NR(i)}^B(x) + \sum_{i=1}^k \pi(i)(x) = 1.$$

If

$$\sum_{i=1}^k \pi_{NR(i)}^B(x) + \sum_{i=1}^k \pi(i)(x) < 1$$

then by definition of  $\pi_{NR}^B$ , we have

$$\sum_{i=1}^k \pi_{NR(i)}^B(x)p_{[i]}(x) + \sum_{i=1}^k \pi(i)(x)p_{[i]}(x) < 1.$$

On the other hand, if

$$\sum_{i=1}^k \pi_{NR(i)}^B(x) + \sum_{i=1}^k \pi(i)(x) > 1$$

then

$$\begin{aligned} \sum_{i=1}^k \pi_{NR(i)}^B(x) + \sum_{i=1}^k \pi(i)(x) &= 1 \\ &+ \sum_{i=1}^k \pi(i)(x)(1 - p_{[i]}(x)). \end{aligned}$$

Therefore, we have

$$r(x, \pi_{NP}^B) \leq r(x, \pi) \text{ for all } \pi \in \mathcal{L}_{NR}^*(\pi, p^*).$$

Corollary 2.2.1. For a given prior  $\pi$  and the loss function

$L = t_1 \mathbb{1}_{\{1\}} + t_2 \mathbb{1}_{\{2\}}$  where  $t_1, t_2 > 0$ , then  $\pi_{NP}^B$  is a non-randomized Bayes

rule wrt the loss function  $l$  for all  $x_1, \dots, x_k$ .

Proof. For the given prior  $\pi$  and the loss function  $l$ , the Bayes risk of any procedure  $\pi$ ,  $\mathcal{C}_{\pi, \text{B}}$ , is, given  $x = (x_1, \dots, x_k)$ ,

$$\begin{aligned} \mathcal{C}_{\pi, \text{B}}(x, l) &= C_1 \sum_{i=1}^k \pi_i(x) + C_2 \sum_{i=1}^k \pi_i(x) \mathbb{P}_{\pi}(x_i \in \mathcal{A}_i) \\ &= C_1 \sum_{i=1}^k \frac{\pi_i(x)}{\pi(x)} + C_2 \sum_{i=1}^k \frac{\pi_i(x)}{\pi(x)} \mathbb{P}_{\pi}(x_i \in \mathcal{A}_i) \\ &= \mathcal{C}_{\pi}(x, \frac{b}{\pi(x)}) \end{aligned}$$

wrt the loss function  $l$ .

Hence  $\mathcal{C}_{\pi, \text{B}}$  is a Bayes-E\* rule wrt the loss function  $l$  for all  $x$ .

### 3.3. Proposed Bayes-E\* Procedure $\pi^*$ in general

Suppose we are interested in the randomized subset selection rule, and we would like to find such a rule which also satisfies the  $E^*$  condition and has the minimum risk wrt the loss function  $l$  for all  $x$  and the prior distribution  $\pi$ .

**Definition 3.3.1.** Given a prior  $\pi$ , we define a Bayes-E\* selection rule,  $\pi^*$ , in which all rules satisfy the condition  $\mathcal{C}_{\pi^*, \text{B}}(x) \leq \mathcal{C}_{\pi, \text{B}}(x)$  for any given observation  $x = (x_1, \dots, x_k)$ , that is,

$$\mathcal{C}_{\pi^*, \text{B}}(x, l) \leq \mathcal{C}_{\pi, \text{B}}(x, l) \text{ for all } x \text{ and } \pi.$$

**Definition 3.3.2.** Given a number  $\delta \in (0, 1)$ , a prior  $\pi$ , a loss function  $l$ , a selection procedure  $\pi^*$  is called  $\delta$ -Bayes-E\* if the procedure  $\pi^*$  is a Bayes rule in the class  $\mathcal{C}_{\pi, \text{B}}$  and

For the sake of convenience, sometimes we will denote the end of  $\mathcal{E}_i(\mathcal{P}^*)$  by

**Definition 2.3.2.** We define a subset selection procedure  $\mathcal{S}$  as follows.

Given a prior  $\pi$  and observation  $x = x_1, \dots, x_n$ ,  $\pi^B$  is defined by

$\beta_1, \beta_2, \dots, \beta_k$

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$$\cdot \frac{B}{(k)}(x) = 1$$

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$$v_{(j)}^B(x) = 1, \quad j \in \bigcup_{i=1}^k p_{(i)j}(x) \cap p^*, \quad j \neq r$$

$$= \begin{cases} \sum_{k=j+1}^K p_{\{i\}}(x) - p^* & \\ \sum_{i=j}^K p_{\{i\}}(x) - p^* & \\ p_{\{i\}}(x) + \sum_{i=j+1}^K p_{\{i\}}(x) - p^* & \end{cases}$$

0 otherwise

Example: If  $E_1 = 1$ ,  $P^* = .90$  and the posterior probabilities are  $p_1 = .95$ ,  $p_2 = .05$ ,  $p_3 = .30$ ,  $p_4 = .15$ , then we select the population  $\pi_1$  with probability  $p_1$ ,  $\pi_2$  with probability  $p_2$ ,  $\pi_3$  with probability  $p_3$ , and we select  $\pi_4$  with probability  $p_4$ , where  $p_4$  is given by

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By Definition 2.3.2 we have

$$\min_{i \in \{1, \dots, k\}} B_{i, \ell^*}(x) \geq \min_{i \in \{1, \dots, k\}} \ell_{\{1\}}^*(x) \ell_{\{1\}}(x) - \ell_{\{1\}}^*(x) \ell_{\{1\}}(x) = 0$$

hence we have the following Lemma.

Lemma 2.3.3.  $\ell_{\{1\}}^*(x) \in \mathcal{C}_{\ell^*}(\ell, P^*)$ .

DEFINITION 2.3.4. we define a subclass  $\mathcal{C}_{\ell^*}^*(\ell, P^*)$  of class  $\mathcal{C}_{\ell^*}(\ell, P^*)$  as

$$\mathcal{C}_{\ell^*}^*(\ell, P^*) = \{ \ell_{\{1\}}^*(x) \ell_{\{1\}}(x) \mid \ell_{\{1\}}(x) \in \mathcal{C}_{\ell^*}(\ell, P^*) \}$$

where  $\ell_{\{1\}}(x) < \dots < \ell_{\{k\}}(x)$  are the ordered  $\ell_i(x)$ 's.

By the definition of  $\mathcal{C}_{\ell^*}^*(\ell, P^*)$  we have the following Lemma.

Lemma 2.3.5.  $\ell_{\{1\}}^*(x) \in \mathcal{C}_{\ell^*}^*(\ell, P^*)$ .

Lemma 2.3.6. for all  $\ell \in \mathcal{C}_{\ell^*}(\ell, P^*)$  there exists  $\ell' \in \mathcal{C}_{\ell^*}^*(\ell, P^*)$  such that  $\ell(x, \ell') = \ell(x, \ell)$  wrt the loss function  $\ell_{\{1\}}$  for all  $x$ .

Theorem 2.3.7. Selection procedure is a Bayesian procedure in  $\mathcal{C}_{\ell^*}(\ell, P^*)$  wrt the loss function  $\ell_{\{1\}}$ .

Proof. Given the observation  $x = x_k$  and any selection procedure

$\ell \in \mathcal{C}_{\ell^*}(\ell, P^*)$  applying  $\ell_{\{1\}}^*(x_k)$  to  $\ell_{\{1\}}(x_k)$  and  $\ell_{\{1\}}^*(x_k) \leq \ell_{\{1\}}(x_k)$  if  $x_k \neq x_1$  hence we have

$$\min_{i \in \{1, \dots, k\}} B_{i, \ell^*}(x_k) \geq 0 \quad \text{for all } x_k.$$

If  $\ell_{\{1\}}(x_k) = 0$ , then we have  $\ell_{\{1\}}^*(x_k) = 0$  and  $\ell_{\{1\}}^*(x_k) \leq \ell_{\{1\}}(x_k)$ .

We will now show that for any  $x_k$

$$\ell_{\{1\}}^*(x_k) \leq \ell_{\{1\}}(x_k)$$

implies

$$P(\mathcal{C}_1^*, x) \geq P(\mathcal{C}_1^B, x) = \pi_{1,1}^*$$

that is

$$\sum_{i=1}^k \pi_{i,1}^*(x) \geq \sum_{i=1}^k \pi_{i,1}^B(x)$$

implies

$$\sum_{i=1}^k \pi_{i,1}^*(x) \rho_{i,1}^*(x) \geq \sum_{i=1}^k \pi_{i,1}^B(x) \rho_{i,1}^B(x)$$

For any  $0 < \alpha < \sum_{i=1}^k \pi_{i,1}^B(x)$ , we have  $0 < \alpha < \pi_{1,1}^*$  when  $\alpha$  is a positive integer and  $0 < \alpha < 1$ .

It is easy to see that the maximum posterior probability of a non-optimal selection of procedure  $i$  with  $\sum_{i=1}^k \pi_{i,1}^*(x) = \alpha$  is

$$\sum_{i=k-\alpha+1}^k \pi_{i,1}^*(x) + \alpha \pi_{k-\alpha,1}^*(x) \geq$$

and it is less than  $\sum_{i=1}^k \pi_{i,1}^B(x) \rho_{i,1}^B(x)$ , since  $\pi_{i,1}^B(x) < \pi_{i,1}^*(x)$ . Therefore  $\pi_{1,1}^*$  is Bayes-P\* procedure in  $\mathcal{C}_1^*$ .

Lemma 2.4. Given the loss function  $L_{\alpha, \beta}$  for all  $\alpha, \beta \in \mathcal{A}$ , then  $\pi_{1,1}^*(x) \rho_{1,1}^*(x)$  is a Bayes-P\* procedure in  $\mathcal{C}_1^*$  if and only if  $\pi_{1,1}^*(x) \rho_{1,1}^*(x) \geq \pi_{1,1}^B(x) \rho_{1,1}^B(x)$  and the posterior risk is with the prior and the loss function  $L_{\alpha, \beta}$ .

Proof. Given  $\pi_{1,1}^*(x) \rho_{1,1}^*(x) \geq \pi_{1,1}^B(x) \rho_{1,1}^B(x)$ .

$$\begin{aligned} & \sum_{i=1}^k \pi_{i,1}^*(x) \rho_{i,1}^*(x) \geq \sum_{i=1}^k \pi_{i,1}^B(x) \rho_{i,1}^B(x) \\ & \text{and } \sum_{i=1}^k \pi_{i,1}^*(x) \rho_{i,1}^*(x) = \pi_{1,1}^* \end{aligned}$$

$$\sum_{i=1}^k \pi_i^{(x)} \nu_i^{(x)}$$

$$\in \mathcal{P}^*$$

hence  $\pi^{(x)} \in \mathcal{P}^*(\pi, \mathcal{P}^*)$ .

Now,

$$\begin{aligned} \pi^{(x, \pi)} &= \sum_{i=1}^k \pi_i^{(x)} \nu_i^{(x)} (1 - \nu_i^{(x)})^{\pi_i^{(x)}} \\ &= \sum_{i=1}^k \pi_i^{(x)} (1 - \nu_i^{(x)})^{1 - \pi_i^{(x)}} \\ &= \sum_{i=1}^k \pi_i^{(x)} (1 - \nu_i^{(x)})^{1 - \nu_i^{(x)}} \\ &= \pi^{(x, \pi)}. \end{aligned}$$

Hence the proof is complete.

**Theorem 2.3.2.** Given the prior  $\pi$  and the observation  $x$ , the procedure  $\pi^{(x, \pi)}$  is a Bayes- $\mathcal{P}^*$  procedure in the class  $\mathcal{P}^*(\pi, \mathcal{P}^*)$  when the prior distribution is  $\pi$ .

**Proof.** By Lemma 2.2.4, it is sufficient to show that

$$\pi^{(x, \pi)} \leq \min_{\pi' \in \mathcal{P}^*} \pi^{(x, \pi')}$$

where

$$\begin{aligned} \pi^{(x, \pi)} &= \sum_{i=1}^k \pi_i^{(x)} (1 - \nu_i^{(x)})^{1 - \nu_i^{(x)}} \\ &= \sum_{i=1}^k \frac{\pi_i^{(x)}}{1 - \nu_i^{(x)}} \frac{1 - \nu_i^{(x)}}{1 - \nu_i^{(x)}} (1 - \nu_i^{(x)})^{1 - \nu_i^{(x)}}. \end{aligned}$$

Let

$$i_0 = \min_{1 \leq i \leq k} \left( \frac{1}{i} \right)^{\lambda^k} = 1$$

$$h_1 = \left( \frac{1}{i_1} \right)^{\lambda^k} (s) < \left( \frac{1}{i_1} \right)^{\lambda^k} (s_0)$$

$$h_2 = \left( \frac{1}{i_2} \right)^{\lambda^k} (s) < \left( \frac{1}{i_2} \right)^{\lambda^k} (s_0),$$

then  $A_1 + i_0 = 1$ .

And we have

$$d_1 = \max \{ h_1 - i_0, -1 \}, \quad \text{if } \left( \frac{1}{i_1} \right)^{\lambda^k} (s_0) < -1,$$

$$d_0 = s_0, \quad \text{if } \left( \frac{1}{i_1} \right)^{\lambda^k} (s_0) \geq -1,$$

$$d_2 = \min \{ h_2 - i_0, -1 \}, \quad \text{if } \left( \frac{1}{i_2} \right)^{\lambda^k} (s_0) < -1,$$

$$i_0 + 1, \quad \text{if } \left( \frac{1}{i_2} \right)^{\lambda^k} (s_0) \geq -1,$$

hence  $d_1 = d_2$ .

Therefore, we have

$$\begin{aligned} e(s, s^k) - e(s, s_0^k) &= \frac{1}{k} \left( \frac{1}{i_1} \right)^{\lambda^k} (s) + \frac{1}{k} \left( \frac{1}{i_2} \right)^{\lambda^k} (s) + \dots + \frac{1}{k} \left( \frac{1}{i_k} \right)^{\lambda^k} (s) \\ &\quad - \frac{1}{k} \left( \frac{1}{i_1} \right)^{\lambda^k} (s_0) - \frac{1}{k} \left( \frac{1}{i_2} \right)^{\lambda^k} (s_0) - \dots - \frac{1}{k} \left( \frac{1}{i_k} \right)^{\lambda^k} (s_0) \\ &= \frac{1}{k} \left( \frac{1}{i_1} \right)^{\lambda^k} (s - s_0) + \frac{1}{k} \left( \frac{1}{i_2} \right)^{\lambda^k} (s - s_0) + \dots + \frac{1}{k} \left( \frac{1}{i_k} \right)^{\lambda^k} (s - s_0) \\ &= \frac{1}{k} \left( \frac{1}{i_1} \right)^{\lambda^k} (s - s_0) + \frac{1}{k} \left( \frac{1}{i_2} \right)^{\lambda^k} (s - s_0) + \dots + \frac{1}{k} \left( \frac{1}{i_k} \right)^{\lambda^k} (s - s_0) \end{aligned}$$

$$\sum_{i \in A_1} \frac{c_i^B(x)}{c_i(x)} \frac{b_i}{b_i(x)} \frac{c_i^B(x)}{c_i(x)} \leq \sum_{i \in A_1} \frac{c_i^B(x)}{c_i(x)} \frac{b_i}{b_i(x)}.$$

$$\star \left( 1 + \frac{c_i^B(x)}{c_i(x)} \right)$$

$$\sum_{i \in A_1} \frac{c_i^B(x)}{c_i(x)} \leq \frac{B}{c_1(x)} (1 + \rho_{A_1, B}).$$

□ by Theorem 2.3.1.

**Corollary 2.3.4.** Procedure  $\star$  is a Bayes- $\star$  rule in  $\mathcal{A}_1$  for the loss function  $l = c_1^B + c_2 l_2$ ,  $c_1, c_2 > 0$ .

**Proof.** Similar to corollary 2.1.1, hence omitted.

#### 2.4. Properties of $\frac{B}{c_1(x)}$ and $\frac{B}{c_2(x)}$

In this section we discuss some properties of the functions  $\frac{B}{c_1(x)}$  and  $\frac{B}{c_2(x)}$ . The following definition of the ordering of distributions was introduced by Lehmann (1953) and further discussed by Lehmann and Bain (1975).

**Definition 2.4.1.** A subset  $A \subseteq \mathbb{R}^d$  is a function of a distribution  $\pi$  if  $\pi_A \leq \pi_B$  for all  $A \in \mathcal{A}$ ,  $\pi_A$  denotes the distribution of  $A$ .

**Definition 2.4.2.** A family of probability distributions  $\{\pi_A\}_{A \in \mathcal{A}}$  is called a  $\star$ -family if  $\pi_A \leq \pi_B$  for all  $A \in \mathcal{A}$ ,  $B \in \mathcal{A}$  and  $\pi_A \leq \pi_B$  for all  $A \in \mathcal{A}$ ,  $B \in \mathcal{A}$ .

$$\pi_A \leq \pi_B \Leftrightarrow \int_A d\pi_A \leq \int_B d\pi_B$$

For all important  $\star$ -families

Let  $f(\cdot, \cdot, \cdot)$  be the p.d.f. of population  $\gamma_1$ . Let  $\pi_{ij}$  be the prior where  $\gamma_{ij}$ 's are mutually independent. Suppose for  $x \in \mathcal{X}$ , we have absolutely continuous posterior c.d.f.  $q_i(\cdot, x)$ . Hence we can write the p.d.f. as

$$q(\cdot, x) = \prod_{j=1}^k q_j(\cdot, x) = \prod_{j=1}^k q_j(\gamma_{ij}, x),$$

Let  $q_i(\cdot, x_i)$  be the posterior c.d.f. associated with  $\gamma_{ij} \sim \pi_{ij}$ ,  $j = 1, \dots, k$ .

**Definition 7.4.3.** The absolutely continuous posterior c.d.f.,  $q_i(\cdot, x)$ ,  $i = 1, \dots, k$ , have the generalized (strictly) stochastic increasing property (G-SIP) if for any  $i_1, i_2, i_3, i_4, j_1, j_2, k$ ,  $x_i < x_j$ , then

$$q_i(\cdot, x_{i_1}) < q_i(\cdot, x_{i_2})$$

Note that if  $\gamma_{ij} \sim \pi_{ij}$  and  $\gamma_{ij} \sim \pi_{ij}$  for all  $j$ ,  $\pi_{ij} \sim \pi_{ij}$ , then the G-SIP is the usual SIP.

**Definition 7.4.4.** A selection procedure  $\gamma$  is monotone consistent if and only if for every  $x \in \mathbb{R}^k$ ,  $x_i < x_j$  implies  $\gamma_{ij} \in \gamma(x)$  and  $\gamma_{ij}$  is a.s. a.s.  $\gamma_{ij}$  is monotone with the increasing of  $x$  almost surely with probability one.

**Theorem 7.4.1.** If the prior  $\pi$  is such that we have  $\pi_{ij} \sim \pi_{ij}$  for all  $i, j$  independent posterior c.d.f.  $q_i(\cdot, x)$  is absolutely continuous, then for every  $x \in \mathbb{R}^k$ ,  $x_i < x_j$  implies

$$\gamma_{ij} \in \gamma(x)$$

and  $\gamma$  is a selection procedure.  $\gamma$  is called a G-SIP selection procedure.

•  $\Gamma(\mathbf{0}, \mathbf{0})$  •

$$\begin{aligned}
 p_i(x) &= P(\gamma_1 < \gamma_{k+1}^*) = \int_{\gamma_1 < \gamma_{k+1}^*} \Phi_{\gamma_1}^*(t) \, dt = \Phi_{\gamma_1}^*(t_{k+1}) \\
 &= \int_{\gamma_1 < \gamma_{k+1}^*} G_i(t) \, dt = M_i(t_{k+1}) - \Phi_{\gamma_1}^*(t_{k+1}) \\
 &= \int_{\gamma_1 < \gamma_{k+1}^*} G_i(t) \, dt \Phi_{\gamma_1}^*(t_{k+1}) \Phi_{\gamma_1}^*(t_{k+1})^* \\
 &= \int_{\gamma_1 < \gamma_{k+1}^*} G_i(t) \, dt \Phi_{\gamma_1}^*(t_{k+1}) \Phi_{\gamma_1}^*(t_{k+1})^* \\
 &= \int_{\gamma_1 < \gamma_{k+1}^*} \frac{d}{dt} \left[ \int_{\gamma_1 < \gamma_{k+1}^*} G_i(t) \, dt \right] \Phi_{\gamma_1}^*(t_{k+1}) \Phi_{\gamma_1}^*(t_{k+1})^* \, dt \\
 &= \int_{\gamma_1 < \gamma_{k+1}^*} G_i(t) \, dt \Phi_{\gamma_1}^*(t_{k+1}) \Phi_{\gamma_1}^*(t_{k+1})^* \, dt
 \end{aligned}$$

since

$$B_j(x) = \frac{b_j}{j!} x^j - 1^{j+1} \cdot B_{j+1}(x) + 1 \cdot \dots$$

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$$\sum_{k=1}^K \left( \lambda_k \right) = \sum_{k=1}^K \left( \lambda_k - \lambda_{\text{opt}} + \lambda_{\text{opt}} \right) = K \lambda_{\text{opt}} - \sum_{k=1}^K \lambda_{\text{opt}} = 0$$

Therefore, the procedures  $\frac{P}{P}$  and  $\frac{B}{B}$  are preferred to

under CMC assumptions, we can re-label the population as  $\pi^* \equiv \pi$  that

Since  $\alpha_1 = \alpha_2 = \alpha_3$ , hence we have  $\alpha_1 \beta_1 \beta_2 = \alpha_2 \beta_1 \beta_2$  and  $\beta_1 \beta_2 \beta_3 = \beta_1 \beta_2 \alpha_3$ .

<sup>10</sup> See also the discussion of the 'feminist' interpretation of the 'right to life' in the next section.

Definition 7.4.6. A selection rule  $\gamma_1$  is called translation invariant if for all  $x \in \mathbb{R}^k$ , and for all  $c \in \mathbb{R}$

$$\gamma_1(x_1 + c, \dots, x_k + c) = \gamma_1(x_1, \dots, x_k) + 1 - c, \quad \forall c \in \mathbb{R}.$$

Definition 7.4.7. A selection rule  $\gamma_1$  is called scale invariant if for all  $x \in \mathbb{R}^k$ , and for all  $c > 0$

$$\gamma_1(x_1 + c, \dots, x_k + c) = \gamma_1(x_1, \dots, x_k) + 1 - c, \quad \forall c \in \mathbb{R}.$$

Theorem 7.4.2. If the posterior distributions  $\alpha_{p_1}(\cdot | \cdot, \cdot, \cdot, \cdot)$  have the shift property, then both selection procedures  $\gamma_1$  and  $\gamma_2$  are "justifiable".

Proof. It is sufficient to show that

$$p_1(x) = \alpha_{p_1}(x | \cdot, \cdot, \cdot, \cdot) \quad \text{whenever } x_1^0 = x_1$$

$$\text{and } 1 - c < x_2 < x_2^0 + \sqrt{1 - c} \quad \forall c \in \mathbb{R}.$$

Suppose  $x_1^0 = x_1$ ,

$$\alpha_{p_1}(x | \cdot, \cdot, \cdot, \cdot) = \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \sum_{i=2}^k (x_i - \mu_i)^2} \alpha_{p_1}(x_1 | \cdot, \cdot, \cdot, \cdot)$$

$$= \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \sum_{i=2}^k (x_i - \mu_i)^2} \alpha_{p_1}(x_1 | \cdot, \cdot, \cdot, \cdot)$$

$$= \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \sum_{i=2}^k (x_i - \mu_i)^2} \alpha_{p_1}(x_1 | \cdot, \cdot, \cdot, \cdot)$$

$$= \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \sum_{i=2}^k (x_i - \mu_i)^2} \alpha_{p_1}(x_1 | \cdot, \cdot, \cdot, \cdot)$$

$$= \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \sum_{i=2}^k (x_i - \mu_i)^2} \alpha_{p_1}(x_1 | \cdot, \cdot, \cdot, \cdot)$$

$$= \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \sum_{i=2}^k (x_i - \mu_i)^2} \alpha_{p_1}(x_1 | \cdot, \cdot, \cdot, \cdot)$$

and for the other  $x_1$ ,  $x_2, \dots, x_k$  the procedure follows.

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SOME RESULTS IN THE THEORY OF SUBSET SELECTION PROCEDURES. (U)

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Definition 2.4.8. Given a number  $P^* \left( \frac{1}{k} < P^* < 1 \right)$ ,  $X = x$  and a prior  $\tau$ , for any selection procedure  $\psi \in \Delta(\tau, P^*)$  the ratio of the posterior probability  $P(CS|\psi, x)$  and the posterior expected selected size  $E(S|\psi, x)$  is called the posterior-efficiency of  $\psi$  and is denoted by  $EFF(\psi|x)$ .

$$EFF(\psi|x) = \frac{P(CS|\psi, x)}{E(S|\psi, x)}.$$

If  $EFF(\psi|x) \geq EFF(\psi'|x)$  for all  $\psi' \in \Delta$  and all  $x$ , then the selection procedure  $\psi$  is called "posterior most efficient" (PME) selection procedure in  $\Delta(\tau, P^*)$ .

Theorem 2.4.3. The non-randomized posterior- $P^*$  selection procedure  $\psi_{NR}^B$  is the PME selection procedure in  $\Delta_{NR}^B(\tau, P^*) = \Delta_{NR}^B$ , given  $\tau, P^*$ .

Proof. By Lemma 2.2.1, for all  $\psi \in \Delta_{NR}$

$$\exists \psi' \in \Delta_{NR}^B \Rightarrow EFF(\psi'|x) \geq EFF(\psi|x) \quad \forall x,$$

hence it is sufficient to show that:

Given  $\tau(\underline{x}), P^*, \underline{x}$ ,  $EFF(\psi_{NR}^B|x) \geq EFF(\psi'|x)$  for all  $\psi' \in \Delta_{NR}^B(\tau, P^*)$ .

We know that, in  $\Delta_{NR}^B(\tau, P^*)$  hence in  $\Delta_{NR}^B$ ,  $\psi_{NR}^B$  always has minimum selected size, i.e.  $\forall \underline{x}, \sum_{i=1}^k \psi_{NRi}^B(\underline{x}) + c = \sum_{i=1}^k \psi_i^B(\underline{x})$  for some integer  $c$ ,  $0 \leq c \leq k - 1$ .

$$\begin{aligned} EFF(\psi'|x) &= \frac{\sum_{i=1}^k \psi_i^B(x) p_{[i]}(x)}{\sum_{i=1}^k \psi_i^B(x)} \\ &< \frac{\sum_{i=1}^k \psi_{NRi}^B(x) p_{[i]}(x) + p_{[k-s-c+1]}(x) + \dots + p_{[k-s]}(x)}{\sum_{i=1}^k \psi_i^B + c} \end{aligned}$$

if  $\psi_{NR}^B(x) = (0, \dots, \underbrace{0, 1, \dots, 1}_{s \text{ terms}})$ .

$$\begin{aligned} \text{EFF}(\psi' | \underline{x}) &\leq \frac{\sum_{i=k-s+1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x}) + c p_{[k-s]}(\underline{x})}{\sum_{i=1}^k \psi_{NR(i)}^B(\underline{x}) + c} \\ &\leq \frac{\sum_{i=1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x})}{\sum_{i=1}^k \psi_{NR(i)}^B(\underline{x})} \\ &= \text{EFF}(\psi_{NR}^B | \underline{x}). \end{aligned}$$

The last inequality is obtained by

$$\begin{aligned} \sum_{i=1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x}) &= \sum_{i=k-s+1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x}) \\ &\geq \left( \sum_{i=k-s+1}^k \psi_{NR(i)}^B(\underline{x}) \right) p_{[k-s]}(\underline{x}). \end{aligned}$$

Theorem 2.4.4. The randomized selection procedure  $\psi^B$  is the PME procedure in  $\mathcal{A}(\tau, P^*) = \mathcal{A}$  for given  $\tau, P^*$ .

Proof. It suffices to show that, given  $\tau, P^*, \underline{x} = \underline{x}$ ,

$$\text{EFF}(\psi^B | \underline{x}) \geq \text{EFF}(\psi' | \underline{x}), \quad \forall \psi' \in \mathcal{A}'.$$

Suppose  $\psi^B(\underline{x}) = (0, \dots, \underbrace{0, 1, \dots, 1}_{s \text{ terms}}) \quad 0 \leq v < 1, 1 \leq s \leq k-1$ .

By theorem 2.3.1 there exists  $c > 0$  such that

$$\sum_{i=1}^k \psi(i)(\underline{x}) + c = \sum_{i=1}^k \psi'(i)(\underline{x}).$$

If  $0 \leq c < 1$ , then

$$\begin{aligned}
 \text{EFF}(\psi' | x) &= \frac{\sum_{i=1}^k \psi'_i(x) p_{[i]}(x)}{\sum_{i=1}^k \psi'_i(x)} \\
 &\leq \frac{\sum_{i=1}^k \psi^B_i(x) p_{[i]}(x) + cp_{[k-s]}(x)}{\sum_{i=1}^k \psi^B_i(x) + c} \\
 &\leq \frac{\sum_{i=1}^k \psi^B_i(x) p_{[i]}(x)}{\sum_{i=1}^k \psi^B_i(x)} \\
 &= \text{EFF}(\psi^B | x).
 \end{aligned}$$

If  $1 \leq c = v' + t + (1-v)$ ,  $t > 0$  integer,  $0 < v' < 1$  then

$$\begin{aligned}
 \sum_{i=1}^k \psi'_i(x) p_{[i]}(x) &= \sum_{i=k-s+1}^k \psi^B_i(x) p_{[i]}(x) + v' p_{[k-s-t+1]}(x) \\
 &\quad + p_{[k-s-t]}(x) + \dots + (1-v) p_{[k-s]}(x) \\
 &\leq \sum_{i=1}^k \psi^B_i(x) p_{[i]}(x) + cp_{[k-s]}(x)
 \end{aligned}$$

hence by the same argument as above we have

$$\text{EFF}(\psi' | x) \leq \text{EFF}(\psi^B | x).$$

Since  $x$  is arbitrary, the result holds for all  $x$ .

## 2.5. Applications to Normal Model

Suppose we have  $k$  populations  $\pi_1, \dots, \pi_k$ ; population  $\pi_i$  has distribution  $N(\mu_i, \sigma_i^2)$ , where  $\sigma_i$ 's are known and  $\mu_i$ 's are unknown. Assume that we

have independent observations  $x_{i1}, \dots, x_{in_i}$ ,  $i = 1, \dots, k$ . Let

$$x_i = \frac{1}{n_i} \sum_{j=1}^{n_i} x_{ij} \text{ and let } \underline{x} = (x_1, \dots, x_k).$$

Suppose we are interested in selecting a subset containing the best (the population having the largest mean) under the posterior-P\* condition, wrt some prior  $\tau = \tau(\mu)$ . Then to find a Bayes-P\* selection procedure is equivalent, in some sense, to finding  $p_i(\underline{x})$ , which is the posterior probability of the event  $\{\mu_i \text{ is the best}\}$ , given observations  $\underline{X} = \underline{x}$ , wrt a given prior  $\tau$ , for all  $i = 1, \dots, k$ .

Case I. Assume that we have a common sample size  $n$  and a common known variance  $\sigma^2$ .

Ia. Suppose we have no prior information about the unknown parameters, and use the "non-informative" (Box and Tiao (1973)) or "locally uniform" prior  $p(\mu_i) \propto c$  for each population.

The posterior density function  $g_i$  of  $\mu_i$ , given  $\underline{x}$  is the normal density with mean  $x_i$  and variance  $\sigma^2/n$ , i.e.,

$$g_i(\mu_i | \underline{x}) = \frac{1}{\sqrt{2\pi}} \exp \left( -\frac{n(\mu_i - x_i)^2}{2\sigma^2} \right).$$

Hence

$$\begin{aligned} p_{[i]}(\underline{x}) &= P(\mu_{(i)} = x_{[k]} | \underline{X} = \underline{x}) \\ &= \int_{-\infty}^{\infty} \frac{1}{\sqrt{2\pi}} \exp \left( -\frac{n}{2\sigma^2} (t + \frac{1}{\sqrt{n}} (x_{[i]} - x_{[j]}))^2 \right) dt, \end{aligned}$$

$i = 1, \dots, k.$

Here  $\mu_{(i)}$  is the quantity corresponding to the  $i^{\text{th}}$  largest observation  $x_{[i]}$ .

Ib. If  $\mu_i$ 's are independent and have the identical prior distribution  $N(\theta_0, \sigma_0^2)$  and  $x_i | \mu_i \sim N(\mu_i, \sigma_1^2/n)$ , then it is well known that the posterior density function  $g_i$  of  $\mu_i$ , given  $\underline{x} = \underline{x}$  is

$$g_i(\mu_i | \underline{x}) \sim N(\bar{\theta}_{x_i}, \xi^2) \text{ with SIP property}$$

where

$$\bar{\theta}_{x_i} = \xi^2 (\sigma_0^{-2} \theta_0 + n \sigma_1^{-2} x_i)$$

$$\xi^2 = (\sigma_0^{-2} + n \sigma_1^{-2})^{-1}.$$

Hence

$$p_{[i]}(\underline{x}) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi(t + \xi n \sigma_1^{-2} (x_{[i]} - x_{[j]})) d\phi(t).$$

The last expression for  $p_{[i]}(\underline{x})$  is the same as that for the non-informative prior whenever  $\sigma_0 \rightarrow \infty$ .

Since  $p_{[i]}(\underline{x}) = p_{[i]}(\underline{x} + \underline{b})$  and since the normal distribution has the strictly SIP, it follows that  $\psi^B$  and  $\psi_{NR}^B$  are "just" a.e. and translation-invariant in both case Ia and Ib.

Case II. Variance  $\sigma_i$ 's are known but  $\sigma_i$ 's and  $n_i$ 's are not all equal.

Iia. Using the non-informative prior  $p(\mu_i) = c$ ,  $i=1, \dots, k$ , we have

$$p_{(i)}(\underline{x}) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi(t + \frac{v_{(i)}}{v_{(j)}} + \frac{x_{[i]} - x_{[j]}}{v_{(j)}}) d\phi(t)$$

where  $v_{(i)} = \frac{\sigma_{(i)}}{n_{(i)}}$ ,  $i = 1, \dots, k$ .  $p_{(i)}$ ,  $v_{(i)}$  and  $n_{(i)}$  are corresponding to  $x_{[i]}$  and we have the following theorem.

Theorem 2.5.1.  $p_{(i)}(x)$  is non-decreasing in  $i$ , i.e.,  $p_{(i)}(x) = p_{[i]}(x)$ .

Remark 2.5.1. From the above formula of  $p_{(i)}(x)$ , it is easy to see, increasing the sample size of the non-best populations will increase the probability that the best population to be selected, however, before doing this, we don't know which one is the best one.

In this case  $\psi^B$  and  $\psi_{NR}^B$  are "just" a.e. and translation-invariant.

Case III. Assume that priors are independent but not identical normal distributions, namely,  $\mu_i \sim N(\theta_i, \sigma_{0i}^2)$ , where  $\theta_i$ 's are not all equal; if the conditional distribution of  $X_i$ , given  $\mu_i$ , is  $N(\mu_i, \frac{\sigma_{1i}^2}{n_i})$ , then the posterior density of  $\mu_i$ , given  $X_i = x_i$  is  $g_i(\mu_i | x_i)$ , which is the probability density function of normal distribution  $N(\bar{\theta}_{x_i}, \sigma_{x_i}^2)$  where

$$\bar{\theta}_{x_i} = \frac{\sigma_{0i}^{-2} \theta_i + n_i \sigma_{1i}^{-2} x_i}{\sigma_{0i}^{-2} + \sigma_{1i}^{-2}}$$

$$\sigma_{x_i}^2 = (\sigma_{0i}^{-2} + \sigma_{1i}^{-2})^{-1}.$$

Hence we have

$$p_i(x) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi\left[t + \frac{\sigma_{0j}^{-2}}{\sigma_{1j}^{-2}} (\bar{\theta}_{x_i} - \bar{\theta}_{x_j})\right] d\phi(t).$$

If  $\sigma_{0i} = \sigma_0$ ,  $\sigma_{1i} = \sigma_1$  and  $n_i = n$ ,  $i = 1, \dots, k$ , then

$$\sigma_{x_i}^2 = \sigma^2 = (\sigma_0^{-2} + \sigma_1^{-2})^{-1} \quad i = 1, \dots, k$$

and

$$p_i(x) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi\left[t + \left(\frac{\sigma_0^{-2}}{\sigma_1^{-2}} + \frac{n(x_i - x_j)}{\sigma_0^2}\right)\right] d\phi(t).$$

#### Case IV. The General Normal Model

Here we consider a more general prior. Suppose we have  $k$  populations, common sample size  $n$  for each population, and common known variance  $\sigma^2 > 0$ . The observation can reduce to  $\underline{X} = (X_1, \dots, X_k)$  where

$$X_i = \sum_{j=1}^n X_{ij}/n, \text{ by sufficiency.}$$

The "Normal Model" is defined as follows:

$$(a) \quad X_i \sim N(\mu, qI), \quad q = \frac{\sigma^2}{n}$$

where  $I$  is the  $k \times k$  identity matrix.

So the  $X$ 's are (conditionally) independent when  $\mu$  is given.

$$(b) \quad \mu \sim N(\theta_0, \gamma I + tU)$$

where  $\theta_0 \in \mathbb{R}$ ,  $\gamma > 0$ ,  $t > -\frac{\gamma}{k}$ ,

$$\mathbf{1} = (1, \dots, 1) \quad \text{and} \quad U = \mathbf{1}' \mathbf{1}.$$

Here  $\gamma > 0$  and  $t > -\frac{\gamma}{k}$  are necessary and sufficient for  $\gamma I + tU$  to be positive definite. This model was chosen by Chernoff and Yahav (1977) ( $t > 0$ ), Gupta and Hsu (1978) and Miescke (1979).

By (a) and (b) we get the posterior distribution of  $\mu$ , given  $\underline{X} = \underline{x}$ , and the distribution of  $X$  as follows:

$$\mu' \underline{x} \sim N(\eta, aI + bU)$$

where

$$\eta = \gamma(q+\gamma)^{-1} \underline{x} + qt((q+\gamma)(q+\gamma+kt))^{-1} \underline{x}' U + q(q+\gamma+kt)^{-1} \mathbf{1}$$

$$a = \gamma q(q+r)^{-1}$$

$$b = q^2 t(q+\gamma)^{-1} (q+\gamma+kt)^{-1}$$

$$\underline{x} \sim N(\underline{m}, (q+\gamma)I + tU)$$

Lemma 2.5.1. Let  $\underline{Y} \sim N(\underline{\mu} + \rho \underline{1}, aI + bU)$  with  $\underline{\mu} \in \mathbb{R}^k$ ,  $\rho \in \mathbb{R}$ ,  $a > 0$  and  $b > -a/k$ . Then there exists a random vector  $\underline{Z} \sim N(\underline{\mu}, aI)$  such that  $h(Y) = h(Z)$  everywhere for every translation-invariant  $h: \mathbb{R}^k \rightarrow \mathbb{R}^k$ .

Proof. (See Miescke (1979)).

With this lemma, it is easy to get

$$p_i(x) = P(\mu_i = \mu_k | \underline{x})$$

$$= \int I_{\{\mu_i = \mu_k\}} d\psi \left( \left( \frac{Y}{q+\gamma} \right) x, \frac{\gamma q}{q+\gamma} I \right)^{(\mu)}$$

where  $\psi(\mu, V)$  is the normal distribution with mean  $\mu$  and variance-covariance matrix  $V$ .

We can rewrite  $p_i(x)$  as

$$p_i(x) = \int_{-\infty}^{\infty} \frac{1}{\sqrt{2\pi}} \phi\left(t + \left(\frac{\gamma}{q(q+\gamma)}\right)^{\frac{1}{2}}(x_i - x_j)\right) d\psi(t).$$

Let  $\gamma = \sigma_0^2$ ,  $q = \sigma^2/n$ , we have

$$p_i(x) = \int_{-\infty}^{\infty} \frac{1}{\sqrt{2\pi}} \phi\left(t + \left(\frac{\sigma_0^2}{\frac{2}{n}(\frac{\sigma^2}{n} + \sigma_0^2)}\right)^{\frac{1}{2}}(x_i - x_j)\right) d\psi(t).$$

The above expression is exactly the same as that of the independent prior Case I, Ib.

Case V. Under normal assumption as before, but suppose  $\sigma_i$ 's are unknown and that neither  $\sigma_i$ 's nor  $n_i$ 's are all equal.

Suppose we have no prior information about  $(\mu, \sigma)$ , for each individual population  $\tau_i$  assign prior  $p(\mu_i, \sigma_i) \cdot \sigma_i^{-1}$  then we have (See Box and Tiao (1973)) that the posterior density of  $\mu_i$ , given  $x_i = x_i^+ (x_{i1}, \dots, x_{in_i})$  is

$$p(\mu_i | x_i) = \frac{(s_i / \sqrt{n_i})^{-1}}{B(\frac{1}{2}v_i, \frac{1}{2})\sqrt{v_i}} [1 + \frac{n_i(\mu_i - x_i)^2}{v_i s_i^2}]^{-\frac{1}{2}(v_i + 1)}$$

where  $s_i^2$  is the sample variance,  $B(\cdot, \cdot)$  is a Beta function and  $v_i = n_i - 1$ .

Hence

$$p(t_i = \frac{\mu_i - x_i}{s_i / \sqrt{n_i}} | x_i) = \frac{1}{B(\frac{1}{2}v_i, \frac{1}{2})\sqrt{v_i}} (1 + \frac{t_i^2}{v_i})^{-\frac{1}{2}(v_i + 1)},$$

which is the density of the student's t distribution with  $v_i (n_i - 1)$  degrees of freedom.

Using this result we can write the formula of  $p_i(x)$  by

$$\begin{aligned} p_i(x) &= p(\mu_i > u_j, \forall j \neq i | x) \\ &= \int_{j \neq i} T v_j \left( t \frac{s_i / \sqrt{n_i}}{s_j / \sqrt{n_j}} + \frac{x_i - x_j}{s_j / \sqrt{n_j}} \right) d T v_i(t) \end{aligned}$$

where  $v_i = n_i - 1$ ,  $i = 1, \dots, k$

$$s_i^2 = \sum_{r=1}^{n_i} (x_{ir} - x_i)^2$$

$T v_i$  is the c.d.f. of t distribution with  $v_i$  degrees of freedom.

When  $v_i$ 's are large, t distribution approaches normal distribution, hence, for large  $n_i$ ,  $i = 1, \dots, k$ , we can replace T by  $\Phi$ .

Case VI. Suppose we are interested in finding a subset which contains the population with the smallest variance; i.e., we define the best population as the one with the smallest variance, and suppose that we have no prior information about  $\sigma$ . In this case, it is reasonable to assume that

$$p(\mu, \sigma) \propto \sigma^{-1}, \text{ if } \mu \text{ is unknown}$$

$$p(\mu, \sigma) \propto 1, \text{ if } \mu \text{ is known.}$$

Let

$$v_j = n_j, v_j s_j^2 = \sum_{r=1}^k (x_{jr} - \mu)^2 \text{ if } \mu \text{ is known}$$

$$v_j = n_j - 1, v_j s_j^2 = \sum_{r=1}^k (x_{jr} - \bar{x}_j)^2 \text{ if } \mu \text{ is unknown, } n_j \geq 1 \quad j = 1, 2, \dots, k$$

$$s^2 = (s_1^2, \dots, s_k^2), \bar{x} = (\bar{x}_{11}, \dots, \bar{x}_{1n_1}, \dots, \bar{x}_{kn_k})$$

and  $Y_{ij}$  be the random variable with c.d.f.  $\chi_{v_j}^2$  which is the  $\chi^2$  distribution with  $v_j$  degrees of freedom.

Then for either case ( $\mu$  known or unknown), we have

$$\begin{aligned} p_i(x) &= P(\sigma_i^2 = s_{[1]}^2 | \bar{x} = x) \\ &= P(\sigma_i^2 \leq v_j^2, \forall j \neq i | \bar{x} = x) \\ &= P\left(\frac{v_j s_j^2}{v_j^2} \leq \frac{v_i s_i^2}{v_i^2} \left(\frac{v_j s_j^2}{v_i^2}\right), \forall j \neq i | s^2 = s^2\right) \\ &= P\left(Y_{ij} \leq \frac{v_i s_i^2}{v_i^2} \left(\frac{v_j s_j^2}{v_i^2}\right), \forall j \neq i | s^2 = s^2\right) \\ &= \int_0^{\infty} \frac{v_i^2}{v_j^2} \frac{v_i^2}{v_i^2} \left(\frac{v_j s_j^2}{v_i^2}\right) d\chi_{v_i}^2(u) \end{aligned}$$

$$= \int_0^{\infty} \int_{\mathbb{R}^k} \psi_j^2(u) \frac{s_j^2}{s_i^2} d\psi_j^2(u) \quad \text{if } n_1 = \dots = n_k = 1.$$

With these  $p_1(x), \dots, p_k(x)$  we can apply Bayes-P\* rules  $\psi^B$  and  $\psi_{NR}^B$  easily.

Lemma 2.5.2. In Case VI,  $\psi^B$  and  $\psi_{NR}^B$  are just \* a.e. and (scale) translation invariant.

\* Here the definition of the "just" property for a selection rule is

$$\psi_i(s_i^2) < \psi_j(s_j^2) \text{ if } s_i^2 > s_j^2, s_j^2 < s_i^2, \forall j \neq i.$$

## 2.6. Comparison of Selection Rules $\psi^B$ and $\psi^M$ in the Normal Location

### Parameter Case

We have  $k$  normal populations with a common known variance  $\sigma^2$  and common sample size  $n$ . For this case Gupta (1956) proposed and studied the procedure  $\psi^M$ .

$\psi^M$ : Select  $x_i$  iff  $x_i > x_{[k]} - d \frac{\sigma}{\sqrt{n}}$   $i = 1, \dots, k$  where  $d = d(\psi^M, \sigma)$  is to be determined by

$$\inf_{\psi \in \mathcal{L}^M} P(CS|\psi^M) = p^*$$

and  $\mathcal{L}$  is the parameter space.

We will show that  $\psi^M \in \mathcal{L}_{NR}(\cdot, p^*)$  where  $\cdot$  is the locally uniform prior distribution. For fixed  $p^*$  and  $k$ , let  $d$  be determined by

$$\int_{-\infty}^{\infty} \psi^{k-1}(t + d) d\psi(t) = p^*. \quad (1.6.1)$$

Let

$\mathcal{X}_i$  = (all possible observed values) =  $\mathbb{R}^k$

$$\mathcal{X}_1 = \{x \in \mathcal{X} | x_{[k]} = d \frac{1}{\sqrt{n}} < x_{[1]}\}$$

$$\mathcal{X}_i = \{x \in \mathcal{X} | x_{[i-1]} < x_{[k]} = d \frac{1}{\sqrt{n}} < x_{[i]}\}, \quad 2 \leq i \leq k$$

$$\mathcal{X}_i^{(1)} = \{x \in \mathcal{X} | x_{[1]} = x_{[i-1]} < x_{[k]} = d \frac{1}{\sqrt{n}} < x_{[i]}\} \subset \mathcal{X}_i$$

$$\mathcal{X}_i^{(2)} = \{x \in \mathcal{X} | x_{[1]} = x_{[i-1]} < x_{[k]} = d \frac{1}{\sqrt{n}} = x_{[i]} = x_{[k-1]}\} \subset \mathcal{X}_i^{(1)}$$

then we have the following theorem.

Theorem 2.6.1. Given a number  $P^* \left( \frac{1}{k} < P^* < 1 \right)$  and locally uniform prior for each population  $\pi_i$ ,  $x = x \in \mathcal{X}_i$ , then

$$P(CS_{i,i}^M, x = x) \geq q^*(i)$$

where

$$q^*(i) = \frac{k-i}{k-1} (1 - P^*) + P^*.$$

Hence

$$x \in \mathcal{X}_{NR}(i, P^*) .$$

Proof. It is sufficient to show that

$$\inf_{x \in \mathcal{X}_i} \sum_{j=1}^k p_{ij,j}(x) = q^*(i) = \frac{k-i}{k-1} (1 - P^*) + P^*.$$

Since  $x \in \mathcal{X}_i$ ,

$$P(CS_{i,i}^M, x) = \inf_{x \in \mathcal{X}_i} P(CS_{i,i}^M, x)$$

$$\inf_{x \in \mathcal{X}_i} \sum_{j=1}^k p_{ij,j}(x) .$$

Without loss of generality we can assume  $\frac{c}{\sqrt{n}} = 1$ .

Since

$$\sum_{i=1}^k p_{[i]}(x) = 1 \quad \forall x \in \mathcal{Z}, \text{ and } \forall i < j$$

$p_{[i]}(x)$  is nonincreasing for all  $x_{[j]}, j \leq i - 1$ , we have

$$\begin{aligned} \inf_{x \in \mathcal{Z}_i} \sum_{i=1}^k p_{[i]}(x) &= \inf_{x \in \mathcal{Z}_i} (1) \sum_{i=1}^k p_{[i]}(x) \\ &= 1 - \sup_{x \in \mathcal{Z}_i} (1) \sum_{j=1}^{i-1} p_{[j]}(x) \\ &= 1 - \sup_{x \in \mathcal{Z}_i} (1) \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \sum_{\substack{t \in \mathcal{Z} \\ j \neq i}} \psi(t + x_{[i]} - x_{[j]}) d\psi(t) \\ &= 1 - \sup_{x \in \mathcal{Z}_i} (1) \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \sum_{\substack{t \in \mathcal{Z} \\ j \neq i}} \psi(t + x_{[i]} - x_{[j]}) \\ &\quad \cdot \sum_{\substack{t \in \mathcal{Z} \\ j \neq i}} \psi(t + x_{[i]} - x_{[j]}) d\psi(t) \\ &= 1 - \sup_{x \in \mathcal{Z}_i} (1) \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \sum_{\substack{t \in \mathcal{Z} \\ j \neq i}} \psi(t + x_{[i]} - x_{[j]}) \\ &\quad \cdot \psi^{i-2}(t) d\psi(t) \\ &= 1 - \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \psi(t - d) \psi^{i-2}(t) d\psi(t) \quad (2.6.7) \end{aligned}$$

$$\begin{aligned}
&= 1 - (i - 1) \int_{-\infty}^{\omega} \phi(t-d) \psi^{k-2}(t) d\phi(t) \\
&= (k - i) \int_{-\infty}^{\omega} \phi^{k-2}(t) \phi(t - d) d\phi(t) \\
&\quad + \int_{-\infty}^{\omega} \phi^{k-1}(t + d) d\phi(t) \quad (2.6.3)
\end{aligned}$$

The superimum of (2.6.3) occurs when  $x \in \mathcal{X}_i^{(2)}$ . The last equality follows from the identity

$$\begin{aligned}
&(k - i) \int \phi^{k-2}(t) \phi(t - d) d\phi(t) \\
&= 1 - \int \phi^{k-1}(t + d) d\phi(t),
\end{aligned}$$

which can be shown by the integration by parts. By (2.6.1), the second term of (2.6.3) equals  $P^*$ ; then use the integration by parts to the first term of (2.6.2), we get

$$\begin{aligned}
\inf_{x \in \mathcal{X}_i} \sum_{j=i}^k p_{[j]}(x) &= \frac{k-i}{k-1} [1 - P^*] + P^* \quad (2.6.4) \\
&= q^*(i).
\end{aligned}$$

Remark 2.6.1. If the procedure  $\psi^M$  selects  $x_{(k)}$  only, i.e.  $X = x \in \mathcal{X}_k$ , then by Theorem 2.6.1 we have  $p_{[k]}(x) \geq P^*$  so that  $\psi^B$  or  $\psi_{NR}^B$  selects  $x_{(k)}$  only. But the converse is not necessarily true.

Remark 2.6.2. For the case  $k = 2$ ,  $\psi_{NR}^B = \psi^M$  a.e. For any given  $X = x$ : if  $x \in \mathcal{X}_2$ , then  $p_{[2]}(x) \geq P^*$ , hence  $\psi^M$  and  $\psi_{NR}^B$  select the population  $x_{(2)}$  associated  $x_{[2]}$ . If  $x \in \mathcal{X}_1$ , and  $x_{[2]} - d \frac{\sigma}{\sqrt{n}} < x_{[1]}$  then  $\psi^M$  and

$\psi_{NR}^B$  select both populations  $\pi_1$  and  $\pi_2$ . Since

$$P(X_{[2]} - d \frac{\pi_1}{\sqrt{n}} = X_{[1]}) = 0,$$

we have  $\psi_{NR}^B = \psi^M$  a.e. .

Remark 2.6.3. The above Theorem and Remark 2.6.1 gives us a lower

bound on the value of  $\sum_{i=1}^k P_{[i]}(x)$ , over all  $x \in \mathbb{C}_i$ . The exact value of the difference of the selected sizes between  $\psi^M$  and  $\psi^B$  depends on the observations.

2.7. Applications to Select  $\max_{1 \leq i \leq k} \pi_i$ ,  $\pi_i = \frac{\mu_i - a}{\sigma_i}$  for Normal

Distribution  $N(\mu_i, \sigma_i^2)$ ,  $i = 1, \dots, k$

Let  $\pi_1, \dots, \pi_k$  be  $k$  independent normal populations with mean  $\mu_i$  and variance  $\sigma_i^2$ , both  $\mu_i$  and  $\sigma_i$  are unknown. For the goal of finding a

random subset which contains the population with maximum  $\pi_i$

for some given constant  $a$ , we assume that apriori  $(\mu_i, \sigma_i)$ ,  $i = 1, \dots, k$  are independent. Suppose we have  $n_i$  independent observations

$x_{i1}, \dots, x_{in_i}$  from  $\pi_i$ , and let  $\bar{x}_i$  be their sample mean,  $i = 1, \dots, k$ .

Let  $y_1, \dots, y_n$  be i.i.d.  $\sim N(\mu, \sigma^2)$ . If no prior information is available to  $(\mu, \sigma)$ , we could assign a locally uniform prior  $p(\mu, \sigma) \propto \sigma^{-1}$  to  $(\mu, \sigma)$ , (see Box and Tiao (1973)). And the posterior joint distribution of  $\mu' = \mu - a$  and  $\sigma$ , given observations

$\mathbf{y} = \mathbf{y}' = (y_1, \dots, y_n)$  is given by

$$P(\mu', \sigma^2 | y) = k \sigma^{-(n+1)} \exp \left\{ -\frac{1}{2\sigma^2} [vs^2 + n(y' - \bar{y})^2] \right\}$$

where

$$\left. \begin{aligned} y' &= y - a, \quad \bar{y} = \sum_{i=1}^n y_i/n \\ vs^2 &= \sum (y_i - \bar{y})^2, \quad v = n - 1 \\ k &= \sqrt{\frac{n}{2\pi}} \left[ \frac{1}{2} \Gamma\left(\frac{v}{2}\right) \right]^{-1} \left( \frac{vs^2}{2} \right)^{\frac{v}{2}}. \end{aligned} \right\} \quad (2.7.1)$$

Let  $t = \sqrt{n}(\mu - a)/\sigma$ , with (2.7.1) the posterior distribution of  $t$ , given  $Y = y$  is

$$p(t | Y = y) = p(t | t)$$

$$= \left\{ 2^{\frac{v}{2}-1} \Gamma\left(\frac{v}{2}\right) \right\}^{-1} \left( \frac{v}{v+t^2} \right)^{\frac{v}{2}} \exp \left\{ -\frac{1}{2} \left( \frac{vt}{v+t^2} \right)^2 \right\} \Gamma(v) I_{v-1} \left( \frac{vt}{v+t^2} \right)^{1/2}$$

where

$$t = \sqrt{n}(y - a)/s, \quad v = n - 1$$

$$I_v(x) = \int_0^{\infty} \left( 2^{\frac{v}{2}} \Gamma\left(\frac{v}{2}\right) \right)^{-1} u^v \exp \left\{ -\frac{1}{2} (u + x)^2 \right\} du.$$

Now, let  $p(\mu_i, \sigma_i) = \sigma_i^{-1}$  be the assigned locally uniform prior to  $(\mu_i, \sigma_i)$ . Then let  $x = (x_{i1}, \dots, x_{1n_1}, \dots, x_{kn_k})$ , we have

$$p_i(x) = P(\mu_i \in \cdot | x)$$

$$= P\left(\frac{\mu_i - a}{\sigma_i} \in \cdot \mid \max_{1 \leq j \leq k} \left( \frac{\mu_j - a}{\sigma_j} \right) \mid x\right)$$

$$\begin{aligned}
 &= P\left(\sqrt{\frac{n_j}{n_i}} \xi_j > \xi_i \mid t\right) \\
 &= \int_{j \neq i} G_{\xi_j} \left( \sqrt{\frac{n_j}{n_i}} z \mid t \right) d G_{\xi_i} (z \mid t) \quad (2.7.1) \\
 &= \int_{j \neq i} G_{\xi_j} (z \mid t) d G_{\xi_i} (z \mid t) \quad \text{if } n_1 = \dots = n_k = n,
 \end{aligned}$$

where  $G_{\xi_i}$  is the posterior c.d.f. of  $\xi_i$  given  $x$  or  $t$ .

By (2.7.2), the Bayes-P\* procedure is completely specified.

If the prior distribution for  $(\mu, \sigma)$  is the conjugate distribution (see Raiffa and Schlaifer (1960)), then

$$\begin{aligned}
 p(\mu, \sigma) &\propto \exp \left\{ -\frac{1}{2\sigma^2} n'(\mu - m')^2 + \frac{1}{\sigma} \right\} \exp \left\{ -\frac{1}{2\sigma^2} (v' + 1) \right\} \\
 &= p(\mu \mid \sigma) p(\sigma)
 \end{aligned}$$

that is

$$p(\mu \mid \sigma) \sim N(m', \sigma^2/n'), \quad n' > 0$$

$$p(\sigma) \sim \frac{v' v''}{2} \sigma^{-2} v', \quad v', v'' > 0.$$

Let

$$x' = \frac{nx + n'm'}{n + n'}, \quad x \text{ is the sample mean.}$$

$$u^2 = (n - 1)s^2 + v'v'' + [(nn')/(n + n')](x - m')^2 / *$$

$$v^* = (n - 1) + v' + 1$$

$$z^* = (n + n')^{1/2}(\mu - a) / *$$

$$t^* = (n + n')^{1/2}(x' - a) / u,$$

the posterior distribution of  $\xi^*$ , given  $x$  is  $p(z^* \mid x) = p(z^* \mid t^*)$  which has the same form as  $p(z \mid t)$ , but replace  $z$ ,  $t$ ,  $\cdot$  by  $z^*$ ,  $t^*$ ,  $\cdot^*$ .

thus, for the conjugate prior case, we get

$$\begin{aligned}
 p_i(\cdot) &= p(\cdot | \mathbf{x}^*, \{t\}^*) \\
 &= \int_{\mathbf{z}^* \in \mathcal{Z}} G_{\mathbf{z}^*}(\mathbf{z}^*) \int_{t^* \in T} z^* t^* dG_{\mathbf{z}^*}(z^*) dG_{t^*}(t^*) \\
 &= \int_{\mathbf{z}^* \in \mathcal{Z}} G_{\mathbf{z}^*}(\mathbf{z}^*) dG_{\mathbf{z}^*}(\mathbf{z}^*) \text{ if } n_i = 1
 \end{aligned}$$

where  $G_{\mathbf{z}^*}$  is the posterior c.d.f. of  $\mathbf{z}^*$  given  $\mathbf{x}$  or  $t^*$ .

Note that (2.7.3) has the same form as (2.7.2), but replace  $\cdot, t$  by  $\cdot^*, t^*$ .

## 2.8. Applications to Poisson Distributions and Poisson Process

### 2.8.1. Poisson Distributions Case

Suppose that  $\gamma_1, \dots, \gamma_k$  are  $k$  independent Poisson populations, where the independent observations  $x_{i1}, \dots, x_{in_i}$  from  $\gamma_i$  have the Poisson density with parameter  $\gamma_i$ ; denoted by  $P(\cdot | \gamma_i)$ ,  $i = 1, \dots, k$ .

Let  $y_1, \dots, y_n$  be i.i.d. with  $p(\cdot | \gamma)$ . If we use non-informative prior  $p(\gamma) \propto \gamma^{-1/2}$  (Box and Tiao (1973)), then given  $\mathbf{y} = (y_1, \dots, y_n)$ , we have the posterior density as follows:

$$p(\gamma | \mathbf{y}) \propto e^{ny - \frac{1}{2}} \exp(-n\gamma)$$

where

$$y = \frac{1}{n} \sum_{i=1}^n y_i \text{ and } \gamma = n(ny + \frac{1}{2}) \Gamma(ny + \frac{1}{2})^{-1}$$

We see that  $2n_i y_i \sim \chi^2_{2n_i+1}$ , the chi-square distribution with  $2n_i+1$  degrees of freedom. Hence by using non-informative prior

$p(\lambda_i) = \lambda_i^{-1/2}$  for each population  $\lambda_i$ , we have

$$\begin{aligned} p_i(x) &= p(\lambda_i = \lambda_{[k]} | x) \\ &= \int_0^\infty \prod_{j \neq i} \chi^2_{n_j}(z) \frac{n_j}{\lambda_j} d\lambda_j \chi^2_{n_i}(z) \end{aligned}$$

where

$$\lambda_i = 2n_i x_i + 1, \quad x_i = \sum_{j=1}^{n_i} x_{ij} / n_i.$$

If  $n_1 = \dots = n_k$ , then

$$p_i(x) = \int_0^\infty \prod_{j \neq i} \chi^2_{n_i}(z) d\lambda_j \chi^2_{n_i}(z).$$

With  $p_i(x)$ ,  $i = 1, \dots, k$ , we can apply Bayes-P\* selection rules

$\lambda^B$  and  $\lambda_{NR}^B$  easily to select a subset which contains the population with the largest parameter  $\lambda$ . On the other hand, if we are interested in selecting the population with the smallest parameter  $\lambda$ , then

$$\begin{aligned} p_i(x) &= \int_0^\infty \prod_{j \neq i} [1 - \chi^2_{n_j}(z)] \frac{n_j}{\lambda_j} d\lambda_j \chi^2_{n_i}(z) \\ &= \int_0^\infty \prod_{j \neq i} [1 - \chi^2_{n_j}(z)] d\lambda_j \chi^2_{n_i}(z) \quad \text{if } n_1 = \dots = n_k. \end{aligned}$$

In this case, the simulation results for selection procedures and  $\lambda_{NR}^B$  are tabulated on Table VII.

### 2.8.2. Poisson Processes Case

Suppose we have  $k$  independent Poisson processes

$\{X^{(1)}(t)\}, \dots, \{X^{(k)}(t)\}$  with expected arrival times equal to

$\lambda_1, \dots, \lambda_k$ , respectively. Hence for the processes  $\{X^{(i)}(t)\}$ , the

probability that there are  $m_i$  arrivals until time  $t_i$  is

$$p(X^{(i)}(t_i) = m_i | \lambda_i, t_i) = \frac{(t_i \lambda_i)^{m_i}}{m_i!} e^{-t_i \lambda_i}$$

$$i = 1, 2, \dots, k, \quad m_i = 0, 1, 2, \dots$$

If there exists no prior information, then we use the non-informative prior  $p(\lambda_i) \propto \lambda_i^{-1/2}$  for all processes. Therefore, we get the posterior density function of  $\lambda_i$ , given  $(m_i, t_i)$  as follows:

$$\begin{aligned} p(\lambda_i | X^{(i)}(t_i) = m_i, t_i) &= p(\lambda_i | m_i, t_i) \\ &= \frac{(t_i \lambda_i)^{m_i + \frac{1}{2} - 1}}{m_i! \left(\frac{1}{2}\right)} e^{-t_i \lambda_i}. \end{aligned}$$

Thus  $2t_i \lambda_i$  has  $\chi^2$  distribution with  $2m_i + 1$  degrees of freedom, given the number  $m_i$  of arrivals before time  $t_i$ .

Let  $m = (m_1, \dots, m_k)$  and  $t = (t_1, \dots, t_k)$ , then it can be shown that the Poisson process  $\{X^{(i)}(t)\}$  has the maximum parameter (or minimum mean waiting time), given  $(m, t)$  is

$$p_i(m, t) = \int_0^\infty \prod_{j \neq i} x_{2m_j+1}^2(y) \frac{t_j}{t_i} d x_{2m_i+1}^2(y) \quad i = 1, \dots, k. \quad (7.2.1)$$

Here we list two special cases which are of interest.

- (a) Observations of all processes are obtained in a common time interval  $[s_i, t + s_i]$ . Since Poisson process is stationary, we can assume that  $s_i = 0$ , and  $t_1 = \dots = t_k = t$ . In this case

$$p_i(m, t) = \int_0^\infty \prod_{j \neq i} x_{2m_j+1}^2(y) d x_{2m_i+1}^2(y)$$

which is independent of  $t$ .

- (b) All  $m_i$ 's are equal, i.e. we fix  $m$  first, then net observations  $t$ .  
Hence

$$p_i(m, t) = \int_0^\infty \prod_{j \neq i} x_{2m+1}^2(y) \frac{t_j}{t_i} d x_{2m+1}^2(y).$$

There is an alternative way to approach the cases (a) and (b). Let  $T_i$  be the waiting time of the  $n$ th arrival in the  $i$ th process, then  $T_i$  has a gamma distribution with density given by

$$p(t) = \frac{\lambda_i^{m_i}}{\Gamma(m_i)} (\lambda_i t)^{m_i-1} e^{-\lambda_i t} \quad t > 0.$$

If we have only non-informative prior  $p(\lambda) \propto \lambda^{-1/2}$ ; then, given  $m_i$  and  $t_i$ ,  $\lambda_i t_i$  has posterior distribution  $\frac{2}{2m_i+1}$ , therefore the formula of  $p_i(m, t)$  we get here is exactly the same as before.

**Remark 2.8.1.** Under non-informative prior, in comparing the subset selection problem in  $k$  Poisson distributions with the problem in  $k$  Poisson processes, it is easily seen that Poisson distributions model is a special case of Poisson processes model, namely,  $t_i = n_i$  where  $n_i$  denotes the sample size of the  $i$ th Poisson population.

### 2.8.3. Relation Between Selection from Poisson Processes and Selection from Populations with Gamma or Exponential Distribution

Suppose we have  $k$  independent populations, the  $i$ th population having the gamma distribution with parameters  $\alpha = m_i$  (known),  $\beta = 1/\lambda_i$  (unknown). Since the random variable  $T_i$ , the waiting time until  $m_i$  arrivals in a Poisson process with parameter  $\lambda_i$ , has a gamma distribution with parameters  $\alpha = m_i$ ,  $\beta = 1/\lambda_i$ . If the  $m_i$ 's are given and if the goals for both selection problems are the same, namely, to select a subset containing the population (process) with the largest parameter  $\beta$ , then it is easily seen that these are identical problems. Note that in the selection problem of Poisson processes,  $m_i$ 's might not be the preassigned values but are given random observations whenever  $t_i$ 's are preassigned values. In this case, the selection problem of Poisson processes is different from that of the gamma distributions.

If the process associated with the minimum parameter  $\beta$  (or the maximum waiting time) is the best, then the posterior probability of process  $x^{(i)}(t)$  to be the best is analogous to the one obtained

before with the modifications that the integrand function

$$\sum_{j \neq i} \frac{2}{x_{2m_j+1}} (y \frac{t_j}{t_i})$$

of (2.8.1) is replaced by

$$\sum_{j \neq i} [1 - \frac{2}{x_{2m_j+1}} (y \frac{t_j}{t_i})].$$

### 2.9. Comparison of the Performance of $\psi^B$ , $\psi_{NR}^B$ , $\psi^M$ and $\psi^{MED}$

Let  $\pi_i$ ,  $i = 1, \dots, k$  be  $k$  independent populations, where  $\pi_i$  has the associated c.d.f.  $F(x, \theta_i) = F(x - \theta_i)$  with unknown location parameter  $\theta_i$ . Let  $f(x, \theta_i) = f(x - \theta_i)$  be the p.d.f. The goal is to find a small (nontrivial) subset which contains the best.

The following subset selection procedure,  $\psi^{MED}$  based on sample medians is due to Gupta and Singh (1980).

$\psi^{MED}$ : Select  $\pi_i$  if and only if  $X_i \leq \bar{X}_{[k]} - d$

where  $\bar{X}_i$  is the median of the  $2m+1$  random observations from population  $\pi_i$  and  $\bar{X}_{[k]} = \max_i \bar{X}_i$ . The value  $d$  is determined by the following equation so that the  $P^*$ -condition is met.

$$\int_{-\infty}^{\infty} G(u + d)^{2m+1} q(u) du = P^*$$

where

$$q(u) = \frac{(2m+1)!}{(m!)^2} [F(u)]^m [1 - F(u)]^m f(u)$$

$$G(u) = I_{F(u)}^{(m+1, m+1)}$$

$I_y(p, q)$  is the incomplete beta function.

In this section we use Monte Carlo simulation techniques to compare the performance of selection procedures  $\psi^B$ ,  $\psi_{NR}^B$ ,  $\psi^M$  and  $\psi^{MED}$  in the normal means problem. Because both rules  $\psi^M$  and  $\psi^{MED}$  are not based on any prior information about the unknown parameters, we assume that the prior distribution  $\pi$  for both  $\psi^B$  and  $\psi_{NR}^B$  is locally uniformly distributed. Since the selection procedure  $\psi^M$  satisfies both the P\*-condition and the posterior-P\* condition wrt the locally uniform priors, it makes sense to compare the Bayes-P\* procedures  $\psi^B$  and  $\psi_{NR}^B$  with  $\psi^M$  and compare  $\psi^M$  with  $\psi^{MED}$  in terms of efficiency which is the ratio of the probability of a correct selection to the expected selected size. For studying the robustness of these four rules,  $\psi^B$ ,  $\psi_{NR}^B$ ,  $\psi^M$  and  $\psi^{MED}$ , we change the true distribution to non-normal distributions, namely, the logistic, Laplace (the double exponential) and the gross error model (the contaminated distribution), but keep the selection procedure unchanged (i.e. still based on the normal assumption). The Monte Carlo simulation results for both equal distances of the parameters and slippage cases are tabulated. In the simulation study all generated random variables are adjusted to have variance 1. Each time we generate five random variables with the given distribution of each population, then apply the selection procedures. The simulation process is repeated 100 times for each random variable. The relative frequency of selecting the population  $\pi_1$  is

used as an approximation to the probability of selecting the population  $\pi_i$ . The sum of relative frequency of selecting each population  $\pi_i$ ,  $i = 1, \dots, k$  is treated as an approximation of the expected selected size. The efficiency  $EFF$  of each selection procedure is approximated by the ratio of relative frequency of selecting the best one to the expected size. The simulation results indicate that in all cases we have the performance

$$\phi^B > \phi_{NP}^B > \phi^M.$$

It should be noted that in the above comparison of the performances, we restrict attention to these rules which satisfy the posterior  $P^*$  condition. For small sample size, the efficiency of rule  $\phi^M$  tends to be larger than  $\phi_{NP}^B$  under  $P^*$ -condition.

Remark 2.9.1. The Laplace distribution has the density function

$$f(x - \mu) = \frac{1}{2} e^{-|x-\mu|} \quad -\infty < x < \infty$$

for which the variance is 2.

The logistic distribution has the density function

$$f(x - \mu) = \frac{e^{-(x-\mu)}}{1 + e^{-(x-\mu)}}^2$$

for which the variance  $\text{Var}(x) = \frac{2}{3}$ .

The gross error model we used has the density function

$$f(x - \mu) = (1 - \beta) \phi(x - \mu + \beta) \cdot \left( \frac{\lambda}{\beta} \right)^{\lambda} \quad \lambda > 0, \beta > 0$$

for which  $\phi$  is the p.d.f. of  $N(0,1)$  and the variance  $\text{Var}(x) = (1 - \beta) + \beta + \beta^2 = 3.25$ .

The efficiency of a selection procedure  $\pi$  is defined by

$$EFF(\pi) = \frac{P_{\pi}(S^*)}{E_{\pi}(S^*)}$$

where  $E_{\pi}(S^*)$  is the expected selected size.

### Discussion and Conclusion

For Table VIII.1 and Table VIII.2 (equal distances case) the  $P^*$  is .99 and .90 respectively, the common sample size  $n = 5$ ,  $k = 5$ . If the  $k$  populations have normal distributions with the unknown parameter configuration  $(\mu_1, \dots, \mu_k, \sigma^2)$ , common variance 1. From both tables the performance based on either the efficiency or the expected selected size is

$$P^* \geq S^* \geq M^*$$

If the posterior- $P^*$  condition is considered, and

$$M^* \geq MED$$

under the  $P^*$ -condition,

when the true distributions are not normal, but the logistic, the Laplace or the trias error model, the results are very close to the normal case, none of the four rules are robust. From Table VIII.3 different  $P^*$ -values than the corresponding ones in Table VIII.1, this is to be expected because the value of  $P^*$  is smaller in the second table.

For Table VIII.1 and VIII.2 (unequal case) the  $P^*$  is .93 and .89 respectively, the common sample size  $n = 5$ ,  $k = 5$ . If the  $k$  populations

have normal distributions with unknown parameter configurations  $(\mu_1, \dots, \mu_k)$ , common variance 1. From both tables the performance based on either the efficiency or the expected selected size is

$$\psi^B \geq \psi_{NR}^B \geq \psi^M$$

if the posterior- $P^*$  condition is considered, and if  $\sqrt{n} = 1$

$$\psi^M > \psi^{MED}$$

under the  $P^*$ -condition.

Note that in both equal distances and slippage cases when  $\psi^M = 1$ , that is the population means are not very close, the procedures  $\psi^B$  and  $\psi_{NR}^B$ , wrt the locally uniform priors, always satisfy not only the posterior- $P^*$  condition but also  $P(\text{CS}|\psi^B \text{ or } \psi_{NR}^B) \geq P^*$ , and the expected selected size of the selection procedure  $\psi^B$  or  $\psi_{NR}^B$  is much less than the selection procedures  $\psi^M$  and  $\psi^{MED}$ . For example, in the normal equal distances case,  $P^* = .99$ ,  $k = 5$ ,  $\sqrt{n} = 4$ ,

$$E(S|\psi^{MED}) - E(S|\psi_{NR}^B) = 0.382;$$

in the normal slippage case,  $P^* = .99$ ,  $k = 5$ ,  $\sqrt{n} = 4$

$$E(S|\psi^{MED}) - E(S|\psi_{NR}^B) = 1.560.$$

TABLE III

For procedures  $\hat{S}$  and  $\hat{S}_2$  and the parameter configurations  $(.5, \dots, 5.0)$  of  $k$  Poisson populations, this table gives the values (based on simulation) of the probability of selecting the population with parameter  $.5i$ ,  $i=1, \dots, 5$ , and the expected selected size  $\hat{S}$ . The prior distribution for each population is  $\text{Dir}(\lambda_1, \dots, \lambda_5)$ .

k	n = 10									
	p*			0.95			0.90			0.75
	$\hat{S}$	NP								
1	.993	1.000	.966	.990	.950	.990	.946	.990	.990	.990
2	.670	.820	.450	.670	.324	.490	.110	.190	.110	.190
5	1.663	1.820	1.415	1.660	1.274	1.480	1.055	1.180	1.055	1.180
7	.968	1.000	.990	1.000	.967	1.000	.929	1.000	.929	1.000
9	.741	.850	.388	.630	.321	.510	.171	.280	.171	.280
3	.205	.330	.120	.130	.065	.110	.021	.050	.021	.050
5	1.944	2.180	1.498	1.760	1.354	1.620	1.121	1.336	1.121	1.336
7	.993	1.060	.981	1.000	.966	.990	.911	.950	.911	.950
9	.745	.829	.560	.730	.259	.400	.175	.290	.175	.290
4	.369	.496	.070	.140	.067	.100	.055	.050	.055	.050
6	.675	.110	.041	.060	.024	.046	.0	.0	.0	.0
8	2.167	2.426	1.653	1.430	1.316	1.530	1.112	1.290	1.112	1.290
1	.996	1.090	.981	.990	.985	1.000	.950	.980	.950	.980
3	.737	.850	.503	.650	.367	.550	.128	.260	.128	.260
5	.355	.479	.102	.160	.133	.210	.013	.050	.013	.050
7	.367	.090	.015	.030	.015	.020	.0	.010	.0	.010
9	.154	.413	.006	.016	.0	.0	.0	.0	.0	.0
2	.154	.413	.1.667	1.667	1.436	1.760	1.397	1.436	1.397	1.436

TABLE VIII. 1

Efficiency (EFF) and Expected Selected Size (ES) (based on simulation) of  $B_M$ ,  $B_{NR}$ ,  $M$  and  $MED$  when the unknown means of the  $k$  populations are  $\beta_1, \dots, \beta_k + (k-1)\delta$ ; the common variance is 1, common sample size  $n = 5$  and the prior for  $B_M$  and  $B_{NR}$  is locally uniformly distributed.

		Normal		Logistic		Tage		Cross error	
		Eff	ES	Eff	ES	Eff	ES	Eff	ES
$n = 3$	B	.254	3.809	.250	3.963	.259	3.772	.248	3.956
$n = 5$	B	.238	4.110	.233	4.290	.238	4.120	.230	4.301
	NR								
	W	.208	4.810	.207	4.840	.210	4.720	.207	4.840
	MED								
	MED	.208	4.810	.202	4.940	.202	4.940	.201	4.930
$n = 7$	B	.333	2.977	.332	3.005	.336	2.941	.329	3.032
	NR								
	W	.250	4.000	.234	4.280	.246	4.030	.233	4.294
	MED								
	MED	.232	4.310	.224	4.460	.217	4.600	.214	4.670
$n = 9$	B	.541	1.847	.541	1.829	.541	1.884	.559	1.779
	NR								
	W	.417	2.400	.417	2.400	.515	2.410	.437	2.290
	MED								
	MED	.369	2.720	.351	2.350	.367	2.760	.369	2.720
$n = 11$	B	.625	1.212	.955	1.162	.621	1.217	.865	1.156
	NR								
	W	.730	1.370	.666	1.240	.746	1.340	.695	1.250
	MED								
	MED	.676	1.420	.694	1.420	.630	1.470	.671	1.470
$n = 13$	B	.777	1.720	.574	1.730	.777	1.770	.671	1.770

TABLE VIII. 2

Efficiency (EFF) and Expected Selected Size (ES) (based on simulation) of  $B$ ,  $B_{NR}$ ,  $M$  and  $MED$  when the unknown means of the  $k$  populations are  $0, \dots, 0 + (k-1)$ ; the common variance is 1, common sample size  $n = 5$  and the prior for  $B$  and  $B_{NR}$  is locally uniformly distributed.

TABLE I

Iteration	Number of nodes	Approximate solution	Lanczos error
1	27	2.000000000000000	0.000000000000000
2	54	1.999999999999999	0.000000000000000
3	108	1.999999999999999	0.000000000000000
4	215	2.000000000000000	0.000000000000000
5	430	1.999999999999999	0.000000000000000
6	855	1.999999999999999	0.000000000000000
7	1710	1.999999999999999	0.000000000000000
8	3420	1.999999999999999	0.000000000000000
9	6840	1.999999999999999	0.000000000000000
10	13680	1.999999999999999	0.000000000000000
11	27360	1.999999999999999	0.000000000000000
12	54720	1.999999999999999	0.000000000000000
13	109440	1.999999999999999	0.000000000000000
14	218880	1.999999999999999	0.000000000000000
15	437760	1.999999999999999	0.000000000000000
16	875520	1.999999999999999	0.000000000000000
17	1751040	1.999999999999999	0.000000000000000
18	3502080	1.999999999999999	0.000000000000000
19	7004160	1.999999999999999	0.000000000000000
20	14008320	1.999999999999999	0.000000000000000
21	28016640	1.999999999999999	0.000000000000000
22	56033280	1.999999999999999	0.000000000000000
23	112066560	1.999999999999999	0.000000000000000
24	224133120	1.999999999999999	0.000000000000000
25	448266240	1.999999999999999	0.000000000000000
26	896532480	1.999999999999999	0.000000000000000
27	1793064960	1.999999999999999	0.000000000000000
28	3586129920	1.999999999999999	0.000000000000000
29	7172259840	1.999999999999999	0.000000000000000
30	14344519680	1.999999999999999	0.000000000000000
31	28689039360	1.999999999999999	0.000000000000000
32	57378078720	1.999999999999999	0.000000000000000
33	114756157440	1.999999999999999	0.000000000000000
34	229512314880	1.999999999999999	0.000000000000000
35	458024629760	1.999999999999999	0.000000000000000
36	916049259520	1.999999999999999	0.000000000000000
37	1832098519040	1.999999999999999	0.000000000000000
38	3664197038080	1.999999999999999	0.000000000000000
39	7328394076160	1.999999999999999	0.000000000000000
40	14656788152320	1.999999999999999	0.000000000000000
41	29313576304640	1.999999999999999	0.000000000000000
42	58627152609280	1.999999999999999	0.000000000000000
43	117254305218560	1.999999999999999	0.000000000000000
44	234508610437120	1.999999999999999	0.000000000000000
45	469017220874240	1.999999999999999	0.000000000000000
46	938034441748480	1.999999999999999	0.000000000000000
47	1876068883496960	1.999999999999999	0.000000000000000
48	3752137766993920	1.999999999999999	0.000000000000000
49	7504275533987840	1.999999999999999	0.000000000000000
50	1500855106797560	1.999999999999999	0.000000000000000
51	3001710213595120	1.999999999999999	0.000000000000000
52	6003420427187560	1.999999999999999	0.000000000000000
53	1200684085437560	1.999999999999999	0.000000000000000
54	240136817087560	1.999999999999999	0.000000000000000
55	48027363417560	1.999999999999999	0.000000000000000
56	96054726835120	1.999999999999999	0.000000000000000
57	192109453670240	1.999999999999999	0.000000000000000
58	384218907340480	1.999999999999999	0.000000000000000
59	768437814680960	1.999999999999999	0.000000000000000
60	1536875629361920	1.999999999999999	0.000000000000000
61	3073751258723840	1.999999999999999	0.000000000000000
62	6147502517447680	1.999999999999999	0.000000000000000
63	12295005034895360	1.999999999999999	0.000000000000000
64	24590010069790720	1.999999999999999	0.000000000000000
65	49180020139581440	1.999999999999999	0.000000000000000
66	98360040279162880	1.999999999999999	0.000000000000000
67	196720080558325760	1.999999999999999	0.000000000000000
68	393440161116651520	1.999999999999999	0.000000000000000
69	786880322233273040	1.999999999999999	0.000000000000000
70	157376064446654680	1.999999999999999	0.000000000000000
71	314752128893309360	1.999999999999999	0.000000000000000
72	629504257786618720	1.999999999999999	0.000000000000000
73	1259008515573237440	1.999999999999999	0.000000000000000
74	2518017031146474880	1.999999999999999	0.000000000000000
75	5036034062292949760	1.999999999999999	0.000000000000000
76	10072068124585899520	1.999999999999999	0.000000000000000
77	20144136249171799040	1.999999999999999	0.000000000000000
78	40288272498343598080	1.999999999999999	0.000000000000000
79	80576544996687196160	1.999999999999999	0.000000000000000
80	161153089993374392320	1.999999999999999	0.000000000000000
81	322306179986748784640	1.999999999999999	0.000000000000000
82	644612359973497569280	1.999999999999999	0.000000000000000
83	1289224719946955138560	1.999999999999999	0.000000000000000
84	2578449439893910277120	1.999999999999999	0.000000000000000
85	5156898879787820554240	1.999999999999999	0.000000000000000
86	10313797759575641108480	1.999999999999999	0.000000000000000
87	20627595519151282216960	1.999999999999999	0.000000000000000
88	41255191038302564433920	1.999999999999999	0.000000000000000
89	82510382076605128867840	1.999999999999999	0.000000000000000
90	165020764153210257755680	1.999999999999999	0.000000000000000
91	330041528306420515511360	1.999999999999999	0.000000000000000
92	660083056612841030752720	1.999999999999999	0.000000000000000
93	1320166113225682061505440	1.999999999999999	0.000000000000000
94	2640332226451364123010880	1.999999999999999	0.000000000000000
95	5280664452902728246021760	1.999999999999999	0.000000000000000
96	10561328905805456492043520	1.999999999999999	0.000000000000000
97	21122657811610912984087040	1.999999999999999	0.000000000000000
98	42245315623221825968174080	1.999999999999999	0.000000000000000
99	84490631246443651936348160	1.999999999999999	0.000000000000000
100	168981262492887303872776320	1.999999999999999	0.000000000000000
101	337962524985774607745552640	1.999999999999999	0.000000000000000
102	675925049971549215491105280	1.999999999999999	0.000000000000000
103	1351850099429848430982210560	1.999999999999999	0.000000000000000
104	2703700198859696861964421120	1.999999999999999	0.000000000000000
105	5407400397719393723928842240	1.999999999999999	0.000000000000000
106	10814800795438787467577684480	1.999999999999999	0.000000000000000
107	21629601590877574935155368960	1.999999999999999	0.000000000000000
108	43259203181755149870310737920	1.999999999999999	0.000000000000000
109	86518406363510299740621475840	1.999999999999999	0.000000000000000
110	173036812727020599481242951680	1.999999999999999	0.000000000000000
111	346073625454041198962485903360	1.999999999999999	0.000000000000000
112	692147250908082397924971806720	1.999999999999999	0.000000000000000
113	1384294501816164795849423613440	1.999999999999999	0.000000000000000
114	2768589003632329591698847226880	1.999999999999999	0.000000000000000
115	5537178007264659183397694453760	1.999999999999999	0.000000000000000
116	1107435601452931836679538896560	1.999999999999999	0.000000000000000
117	2214871202905863673359077793120	1.999999999999999	0.000000000000000
118	4429742405811727346718155586240	1.999999999999999	0.000000000000000
119	8859484811623454693436311172480	1.999999999999999	0.000000000000000
120	1771896962324690938687262234960	1.999999999999999	0.000000000000000
121	3543793924649381877374524469920	1.999999999999999	0.000000000000000
122	7087587849298763754749048939840	1.999999999999999	0.000000000000000
123	14175175698597327509498097879680	1.999999999999999	0.000000000000000
124	28350351397194655018996195759360	1.999999999999999	0.000000000000000
125	56700702794389310037992391518720	1.999999999999999	0.000000000000000
126	113401405988778620075984783037440	1.999999999999999	0.000000000000000
127	226802811977557240151969566074880	1.999999999999999	0.000000000000000
128	453605623955114480303939132149760	1.999999999999999	0.000000000000000
129	907211247910228960607878264299520	1.999999999999999	0.000000000000000
130	181442247820045792121575652859840	1.999999999999999	0.000000000000000
131	362884495640091584243151305719680	1.999999999999999	0.000000000000000
132	725768991280183168486302611439360	1.999999999999999	0.000000000000000
133	145153798256036633697264522287720	1.999999999999999	0.000000000000000
134	290307596512073267394529044575440	1.999999999999999	0.000000000000000
135	580615193024146534789058089150880	1.999999999999999	0.000000000000000
136	1161230386048293069578117178301760	1.999999999999999	0.000000000000000
137	2322460772096586139156234356603520	1.999999999999999	0.000000000000000
138	4644921544193172278312468713207040	1.999999999999999	0.000000000000000
139	9289843088386344556624937426414080	1.999999999999999	0.000000000000000
140	1857968617677268911324985485228160	1.999999999999999	0.000000000000000
141	3715937235354537822649970970456320	1.999999999999999	0.000000000000000
142	7431874470709075645299941940912640	1.999999999999999	0.000000000000000
143	1486374894141815129559883988185280	1.999999999999999	0.000000000000000
144	2972749788283630259119767976370560	1.999999999999999	0.000000000000000
145	5945499576567260518239535952741120	1.999999999999999	0.000000000000000
146	1189099915313452103647857190548240	1.999999999999999	0.000000000000000
147	2378199830626904207295714381096480	1.999999999999999	0.000000000000000
148	4756399661253808414585428762192960	1.999999999999999	0.000000000000000
149	9512799322507616829170857524385920	1.999999999999999	0.000000000000000
150	1902559864501523365834171504877920	1.999999999999999	0.000000000000000

TABLE IX. 1

Estimated average (standardized) selected size (ES) (based on simulation) of  $B_{NR}$ ,  $B_M$  and  $B_{MED}$  when the true variance of the  $k$ th population is  $\sigma_k^2 = 1 + \frac{1}{k}$ ; the common variance is 1, common sample size  $n = 5$  and type prior for  $B$  and  $B_M$  is locally uniformly distributed.

TABLE IV. 1  
 $\alpha = 5$ ,  
 $P^* = .35$

n	k	Normal		Logistic		Laplace		Gross error	
		EFF	ES	EFF	ES	EFF	ES	EFF	ES
2	2	4.132	.218	4.253	.226	4.203	.225	4.361	
3	3	4.527	.212	4.580	.214	4.530	.212	4.670	
4	4	4.373	.202	4.950	.200	4.900	.202	4.890	
5	5	4.420	.202	4.940	.202	4.960	.200	5.000	
6	6	4.167	.240	4.153	.240	4.062	.231	4.264	
7	7	4.270	.224	4.460	.223	4.430	.216	4.580	
8	8	4.367	.203	4.930	.203	4.670	.214	4.900	
9	9	4.397	.201	4.970	.200	4.980	.200	5.000	
10	10	4.546	.274	3.643	.263	3.701	.265	3.381	
11	11	4.972	.253	3.360	.242	4.090	.269	3.720	
12	12	4.626	.211	4.750	.214	4.680	.212	4.510	
13	13	4.287	.207	4.510	.203	4.470	.207	4.270	
14	14	4.346	.205	4.520	.205	4.510	.205	4.350	
15	15	4.346	.204	4.520	.204	4.510	.204	4.350	
16	16	4.346	.204	4.520	.204	4.510	.204	4.350	
17	17	4.346	.204	4.520	.204	4.510	.204	4.350	
18	18	4.346	.204	4.520	.204	4.510	.204	4.350	
19	19	4.346	.204	4.520	.204	4.510	.204	4.350	
20	20	4.346	.204	4.520	.204	4.510	.204	4.350	

TABLE IX. 2

Estimated average and standard selected size (SES) (based on simulation) of  $B$ ,  $B_{\text{M}}$ ,  $N_{\text{H}}$ ,  $N_{\text{M}}$  and  $N_{\text{EO}}$  when error variances and covariances in simulations are  $0.01$ ,  $0.01$ ,  $0.01$ ; the common variance is  $1$ ; common sample size is  $100$ ; and there are  $10$  observations in  $B$  and  $N_{\text{H}}$ ; initially uninformative.

N	M	Logistic	Logistic	Estimate		Gross error	
				EF	ES	EF	ES
10	10	1.250	1.250	.238	2.734	.256	2.856
10	20	1.262	1.262	.228	3.160	.240	3.250
10	30	1.264	1.264	.211	4.400	.212	4.490
10	40	1.264	1.264	.202	4.840	.204	4.900
10	50	1.264	1.264	.194	5.354	.203	5.471
10	60	1.264	1.264	.186	5.905	.205	6.015
10	70	1.264	1.264	.178	6.496	.206	6.616
10	80	1.264	1.264	.170	7.127	.207	7.247
10	90	1.264	1.264	.162	7.797	.208	7.917
10	100	1.264	1.264	.154	8.498	.209	8.618
10	120	1.264	1.264	.146	9.249	.210	9.369
10	140	1.264	1.264	.138	10.040	.211	10.160
10	160	1.264	1.264	.130	10.871	.212	10.991
10	180	1.264	1.264	.122	11.742	.213	11.862
10	200	1.264	1.264	.114	12.653	.214	12.773
10	220	1.264	1.264	.106	13.604	.215	13.724
10	240	1.264	1.264	.098	14.595	.216	14.715
10	260	1.264	1.264	.090	15.516	.217	15.636
10	280	1.264	1.264	.082	16.467	.218	16.587
10	300	1.264	1.264	.074	17.448	.219	17.578
10	320	1.264	1.264	.066	18.459	.220	18.599
10	340	1.264	1.264	.058	19.500	.221	19.630
10	360	1.264	1.264	.050	20.571	.222	20.711
10	380	1.264	1.264	.042	21.672	.223	21.812
10	400	1.264	1.264	.034	22.803	.224	22.983
10	420	1.264	1.264	.026	23.964	.225	24.144
10	440	1.264	1.264	.018	25.155	.226	25.315
10	460	1.264	1.264	.010	26.376	.227	26.546
10	480	1.264	1.264	.002	27.627	.228	27.797
10	500	1.264	1.264	-.006	28.908	.229	29.078
10	520	1.264	1.264	-.014	30.219	.230	30.389
10	540	1.264	1.264	-.022	31.550	.231	31.720
10	560	1.264	1.264	-.030	32.891	.232	33.062
10	580	1.264	1.264	-.038	34.242	.233	34.413
10	600	1.264	1.264	-.046	35.603	.234	35.774
10	620	1.264	1.264	-.054	36.974	.235	37.145
10	640	1.264	1.264	-.062	38.355	.236	38.526
10	660	1.264	1.264	-.070	39.746	.237	40.017
10	680	1.264	1.264	-.078	41.147	.238	41.328
10	700	1.264	1.264	-.086	42.568	.239	42.759
10	720	1.264	1.264	-.094	44.009	.240	44.240
10	740	1.264	1.264	-.102	45.460	.241	45.691
10	760	1.264	1.264	-.110	46.921	.242	47.162
10	780	1.264	1.264	-.118	48.402	.243	48.633
10	800	1.264	1.264	-.126	50.003	.244	50.204
10	820	1.264	1.264	-.134	51.624	.245	51.795
10	840	1.264	1.264	-.142	53.265	.246	53.456
10	860	1.264	1.264	-.150	54.926	.247	55.097
10	880	1.264	1.264	-.158	56.607	.248	56.768
10	900	1.264	1.264	-.166	58.308	.249	58.439
10	920	1.264	1.264	-.174	60.029	.250	60.110
10	940	1.264	1.264	-.182	61.760	.251	61.881
10	960	1.264	1.264	-.190	63.501	.252	63.652
10	980	1.264	1.264	-.198	65.252	.253	65.423
10	1000	1.264	1.264	-.206	67.013	.254	67.194

BIBLIOGRAPHY

## BIBLIOGRAPHY

- Alam, E. (1973). On a multiple decision rule. *Ann. Statist.*, 1, 750-755.
- Andersen, E. S. (1953). On the fluctuation of sum of random variables. *Math. Scand.*, 1, 263-285.
- Ayer, M., Brunk, H. D., Eaton, G. M., Reid, W. T. and Silverman, E. (1955). An empirical distribution function for sampling with incomplete information. *Ann. Math. Statist.*, 26, 641-652.
- Bahadur, R. R. (1956). On the problem in the theory of estimation. *Ann. Math. Statist.*, 27, 362-375.
- Bahadur, R. R. and Robbins, H. (1956). The problem of the adaptive mean. *Ann. Math. Statist.*, 27, 469-487. *Corrigendum*, 28, 149-150.
- Bairamov, R. E., Bartholomew, D. J., Bremner, J. M. and Tsiatis, A. (1973). *Statistical Inference Under Order Restrictions*. John Wiley, New York.
- Barron, A. M. and Gupta, S. S. (1971). A class of non-elliptical sequential multiple decision procedures. *Operations Research Verfahren* (Ed. Bemby, Künnzi and Schubert). Verlag Anton Hain, Meisenheim am Glan, Germany, pp. 11-7.
- Bechhofer, R. E. (1964). A multiple-sample multiple decision procedure for ranking means of normal populations with known variances. *Ann. Math. Statist.*, 35, 166-181.
- Bechhofer, R. E. (1967). A sequential multiple decision procedure for selecting the best mean of several normal populations with common unknown variance, in use with various expected deviations. *Biometrika*, 54, 308-319.
- Bechhofer, R. E., Tamhane, S. C. and Folks, M. (1967). A two-sample multiple-decision procedure for ranking means of normal populations with common unknown variances. *Biometrika*, 54, 179-197.
- Bechhofer, R. E., Tamhane, S. C. and Folks, M. (1968). A two-sample multiple-decision procedure for ranking means of normal populations with common unknown variances. *Biometrika*, 55, 319-330.

- Sechhofer, R. E. and Sobel, M. (1961). A single-sample multiple decision procedure for ranking variances of normal populations (prelim. report). *Ann. Math. Statist.*, 32, 107-113.
- Sherger, R. L. (1971). Minimax, admissible and sigma-minimax multiple decision rules. Ph.D. Thesis, Mimeo. Ser. No. 499, Dept. of Math. Statist., Purdue Univ., West Lafayette, Indiana.
- Sherger, R. L. (1972). Minimax subset relation for two-variance multiple subset size. *Ann. Statist.*, 7, 117-137.
- Sherger, R. L. and Gupta, S. S. (1970). Minimax subset selection rules with application to model variance problems. *Ann. Math. Statist.*, 7, 31-36.
- Sickel, R. J. and Yahav, J. A. (1973). On selecting a subset of a population. *Statistical Decision Theory and Related Inference Problems, I* (Eds. S. S. Gupta and B. N. Prasad). Academic, New York, pp. 311-326.
- Box, G. E. P. and Tiao, G. C. (1973). Bayesian Inference in Statistical Analysis. Addison Wesley, Massachusetts.
- Brostrom, B. (1971). An improved procedure for selecting a subset of a population. *Statistical Decision Theory and Related Inference Problems, I* (Eds. S. S. Gupta and B. N. Prasad). Academic, New York, pp. 311-326.
- Tuckill, R. C. (1971). Asymptotically nonparametric sequential selection procedures. Mimeo. Ser. No. 944, Inst. of Stat. and Math., Univ. of North Carolina, Chapel Hill.
- Thompson, R. and Yahav, J. (1971). A subset selection rule based on evaluating a new criterion. *Statistical Decision Theory and Related Topics, I* (Eds. S. S. Gupta and B. N. Prasad). Academic, New York, pp. 33-110.
- Wells, J. J. (1965). Multiple decision procedures: from an empirical decision approach. Ph.D. Thesis, Mimeo. Ser. No. 451, Dept. of Math. Statist., Purdue Univ., West Lafayette, Indiana.
- Zeely, J. J. and Gupta, S. S. (1969). On the properties of subset selection procedures. *Statist. Sci.*, 4, 39-46.
- Desu, M. M. and Sobel, M. (1960). A fixed subset size approach to a selection problem. *Biometrika*, 47, 401-410. (Corrections and amendments: 48 (1961), 185).
- Desu, M. M. and Sobel, M. (1971). Nonparametric procedures for selecting fixed-size subsets. *Statistical Decision Theory and Related Topics (Eds. S. S. Gupta and B. N. Prasad)*. Academic, New York, pp. 266-276.

- Samuels, L. W. (1951). A multiple comparison procedure for comparing several treatments with a control. *J. Amer. Statist. Assoc.*, 46, 108-113.
- Keller, W. (1971). *An Introduction to Probability Theory and Its Applications*, Vol. II, 2nd Edition. J. Wiley & Sons, New York.
- Robbins, J. D., Barb, T. and Lovel, R. (1972). *Selection and Ranking Populations*. Wiley, New York.
- Sattarzadeh, M. (1967). On selection and ranking procedures in multivariate normal populations. Ph.D. Thesis, Inst. of Statist., Purdue Univ., West Lafayette, Indiana.
- Swaminathan, M. and Gupta, J. S. (1970). Selection Procedures in multivariate normal distributions in terms of measures of distance. *Econometrics*, 37, 105-117.
- Seely, P. and Rubin, R. (1977). On selecting a subset (certainly) the best population - a Bayesian approach. *Ann. Statist.*, 5, 692-702.
- Swaminathan, M. (1956). On a decision rule for a profile in ranking problems. Ph.D. Thesis (Mimeo. Ser. No. 152). Inst. of Statist., Univ. of North Carolina, Chapel Hill.
- Swaminathan, M. (1963). On a selection and ranking procedure for multivariate normal populations. *Ann. Inst. Statist. Math.*, 14, 179-197.
- Swaminathan, M. (1964). On some multiple decision (selection) procedures. *Technometrics*, 6, 335-345.
- Swaminathan, M. (1967). On selection and ranking procedures. *Technometrics*, 9, 151-155.
- Gupta, J. S. and Swaminathan, M. (1971). On unbiased selection and ranking procedures for selecting populations close to a target population. *Technometrics*, 13, 331-341.
- Gupta, J. S. and Swaminathan, M. (1970). On the performance of some unbiased selection procedures. *Technometrics*, 12, 351-361.
- Gupta, J. S. and Swaminathan, M. (1970). On subset selection procedures for Poisson populations and some applications to multiple decision and selection problems. *Applied Statistics* (Ed. R. R. Hinde and J. G. Pollard, Academic, London), pp. 312-323.
- Gupta, J. S. and Swaminathan, M. (1976). On some unbiased selection and ranking procedures in sequential subset selection procedures. *Technometrics*, 18, 371-378. *Inference and Statistical Decision*, Vol. 2 (Ed. M. M. Raghavarao), New York, pp. 101-111.

- Gupta, S. S. and Huang, D. Y. (1976). Selection procedures for the means and variances of normal populations: unequal sample size case. *Sankhyā Ser. B*, 38, 112-128.
- Gupta, S. S. and Kim, W. C. (1980).  $\beta$ -minimax and minimax decision rules for comparison of treatments with a control. To appear in *Recent Developments in Statistical Inference and Data Analysis*, North Holland Publishing Company, 1980, 56-72.
- Gupta, S. S. and McDonald, G. C. (1972). On some class of selection procedures based on ranks. *Nonparametric Techniques in Statistical Inference* (Ed. M. L. Puri), Cambridge University Press, 1972, pp. 491-511.
- Gupta, S. S. and Miescke, K. J. (1978). On subset selection procedures for ranking means of three normal distributions. *Unpublished Ser. No. 78-19*, Dept. of Statist., Purdue Univ., West Lafayette, Indiana. To appear in *Sankhyā Ser. A*.
- Gupta, S. S. and Miescke, K. J. (1979). On the best subset sample size configurations in certain two-stage selection procedures. *Unpublished Ser. No. 79-6*, Dept. of Statist., Purdue Univ., West Lafayette, Indiana.
- Gupta, S. S. and Nagel, K. (1971). On some contributions to multiple decision theory. *Statistical Decision Theory and Related Inference* (Eds. S. S. Gupta and J. N. Keeyell), Academic, New York, 1971.
- Gupta, S. S., Nagel, K. and Panchapakesan, S. (1971). On the use of statistics from equally correlated normal random variables. *Biometrika*, 60, 403-413.
- Gupta, S. S. and Panchapakesan, S. (1973). On a class of subset selection procedures. *Ann. Math. Statist.*, 43, 914-927.
- Gupta, S. S. and Panchapakesan, S. (1979). *Multiple Decision Theory: Theory and Methodology of Selection and Ranking Population*. John Wiley, New York.
- Gupta, S. S. and Singh, A. K. (1980). On rules based on sample size for selection of the largest three population. *Statistica, Teoria, Meth.*, 19(12), 157-179.
- Gupta, S. S. and Sobel, M. (1961). On selecting a subset which contains all populations better than a standard. *Ann. Math. Statist.*, 32, 235-244.
- Gupta, S. S. and Sobel, M. (1962a). On selecting a subset containing the population with the smallest variance. *Biometrika*, 49, 169-178.
- Gupta, S. S. and Sobel, M. (1962b). On the smallest of  $k$  correlated test statistics. *Biometrika*, 49, 369-378.

- Gupta, S. S. and Shaffer, J. B. (1971). On some selection and ranking procedures with applications to multivariate populations. *Proc. in Probability and Statistics*, Ed. R. C. Bose et. al., University of North Carolina Press, Chapel Hill, pp. 397-421.
- Hwang, F. G. (1974). A simple sequentially rejective multiple test procedure. *Biometrika*, 61, 123-128.
- Hwang, P. (1979). Some contributions to gamma-minimax and empirical Bayes selection procedures. Ph.D. thesis (Mimeo. Ser. No. 361), Dept. of Statist., Purdue Univ., West Lafayette, Indiana.
- Hwang, P. Y. (1975). Some contributions to fixed sample and sequential multiple decision (selection and ranking) theory. Ph.D. thesis (Mimeo. Ser. No. 363), Dept. of Statist., Purdue Univ., West Lafayette, Indiana.
- Hwang, W. T. (1972). Some contributions to sequential selection and ranking procedures. Ph.D. thesis (Mimeo. Ser. No. 298), Dept. of Statist., Purdue Univ., West Lafayette, Indiana.
- Kruskal, J. B. (1964). Nonmetric multidimensional scaling: a nonmetric method. *Psychometrika*, 29, 115-129.
- Lehmann, E. L. (1952). Testing multiparameter hypotheses. *Ann. Math. Statist.*, 23, 541-552.
- Lehmann, E. L. (1955). Ordered families of distributions. *Ann. Math. Stat.*, 26, 413-428.
- Lehmann, E. L. (1961). Some Model I problems of selection. *Ann. Math. Statist.*, 32, 990-1012.
- Mahamudulu, D. M. (1967). Some fixed-sample ranking and selection problems. *Ann. Math. Statist.*, 38, 1079-1091.
- McDonald, G. C. (1969). On some distribution-free ranking and selection procedures. Ph.D. thesis (Mimeo. Ser. No. 174), Dept. of Statist., Purdue Univ., West Lafayette, Indiana.
- Mielke, E. D. (1979). Bayesian subset selection for additive and linear loss functions. *Comput. Statist. - Theor. Meth.*, AP(1), 1205-1226.
- Miettinen, T. (1948). A k-sample slippage test for an extreme population. *Ann. Math. Statist.*, 19, 58-65.
- Nand, E. (1970). On subset selection rules with certain optimality properties. Ph.D. thesis (Mimeo. Ser. No. 222), Dept. of Statist., Purdue Univ., West Lafayette, Indiana.
- Nand, E. (1974). On subset selection rules for comparing several normal distributions. *Ann. Statist.*, 4, 529-539.

- Naik, A. D. (1977). Some subset selection problems. *Journal of Multivariate Analysis*, 7, 955-966.
- Naik, A. D. (1978). On selection procedure based on the exceedence probability. *Calcutta Statist. Assoc. Bull.*, 19, 40.
- Paulson, E. (1949). A multiple decision procedure for certain problems in the analysis of variance. *Ann. Math. Statist.*, 20, 96-106.
- Paulson, E. (1962). A sequential procedure for comparing several experimental categories with a standard or control. *Ann. Math. Statist.*, 33, 438-443.
- Paulson, E. (1963). A sequential decision procedure for choosing one of  $k$  hypothesis concerning the unknown mean of a normal distribution. *Ann. Math. Statist.*, 34, 549-554.
- Paulson, E. (1964). A sequential procedure for selecting the  $m$  categories with the largest mean from  $k$  normal populations. *Ann. Math. Statist.*, 35, 174-180.
- Paulson, E. (1967). Sequential procedures for selection of the  $m$  best of several binomial populations. *Ann. Math. Statist.*, 38, 1-14.
- Rabinowitz, P. and Weiss, G. (1960). Tables of abscissae and weights for numerical evaluation of integrals of the form  $\int_0^1 e^{-xt} f(t) dt$ . *Math. Tables and other Aids to Comput.*, Vol. XIII, 65, 155-164.
- Rizvi, M. H. and Sobel, M. (1967). Nonparametric procedures for selecting a subset containing the population with the largest mean. *Ann. Math. Statist.*, 38, 726-741.
- Santner, T. J. (1975). A restricted subset selection approach to ranking and selection problems. *Ann. Statist.*, 3, 334-346.
- Seal, R. C. (1955). On a class of decision procedures for ranking the mean of normal populations. *Ann. Math. Statist.*, 26, 321-332.
- Seal, R. C. (1957). An optimum decision rule for ranking the means of several populations. *Calcutta Statist. Assoc. Bull.*, 1, 11-16.
- Sobel, M. (1967). Nonparametric procedures for selecting the  $m$  best distributions with the largest mean. *Ann. Math. Statist.*, 38, 1604-1616.
- Sobel, M. and Rivot, M. (1967). Selecting the best one of several binomial populations. *Bell System Tech. J.*, 46, 925-936.
- Studden, W. J. (1967). On selecting a subset of  $k$  populations containing the best. *Ann. Math. Statist.*, 38, 1617-1629.

- Tamhane, A. C. and Bechhofer, R. E. (1977). A two-stage minimax procedure with screening for selecting the largest normal mean. *Commun. Statist. - Theor. Meth.*, A6(11), 1003-1033.
- Tamhane, A. C. and Bechhofer, R. E. (1979). A two-stage minimax procedure with screening for selecting the largest mean (I): An improved PCS lower bound and associated tables. *Commun. Statist. - Theor. Meth.*, A8(4), 337-358.
- Tukey, J. W. (1960). A survey of sampling from contaminated distributions. *Contributions to Probability and Statistics* (Eds. J. W. Tukey and others). Stanford Univ. Press, 39, 448-458.

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20. ABSTRACT (list in reverse side if necessary and identify by block number) The problem of selection and ranking (ordering) problems in statistics is often studied to gain the classical truth of homogeneity assumption. Then the problem is to select the treatments where the experimenter is interested in comparing the efficiency of treatments or processes with the goal of selecting the best treatment within a population.		
Chapter 1 of this thesis considers the problem of selecting the best treatment within a population if it is better than a control under the assumption that		

an ordering prior. Here, by an ordering prior we mean that there exists a known simple or partial order relationship among the unknown parameters of the treatments (excluding the control). Three new selection procedures are proposed and studied. These procedures do meet the usual requirement that the probability of a correct selection is greater than or equal to a pre-specified number  $P^*$ . Two of the three procedures use the isotonic regression estimator of sample means of the  $k$  treatments with respect to the given ordering prior. Tables which are necessary to carry out the selection procedure with this approach for the selection of unknown means of normal populations and of non-normal populations are given. Monte Carlo comparisons on the performance of the three procedures for the normal or gamma mean problem were carried out in various selected cases. The results of this study seem to indicate that the procedures based on isotonic estimators always have superior performance, especially when there are more than one bad population. (in comparison with the existing)

Chapter II deals with a new 'Bayes- $P^*$ ' approach about the procedure of selecting a subset which contains the 'best' of  $k$  populations, where by best we mean the (unknown) population with the largest unknown mean. The (non-randomized) Bayes- $P^*$  rule refers to a rule with minimax risk in the class of (non-randomized) rules which satisfy the condition that the probability of the ability of selecting the best is at least equal to  $P^*$ . Given the prior of the unknown parameters, two 'Bayes- $P^*$ ' subset selection procedure

$J^B$  and  $J^B_{NR}$  (randomized and non-randomized, respectively) under quadratic loss functions are obtained and compared with one classical subset selection procedure  $J^M$ . The comparisons of the performance of  $J^B$  with  $J^B_{NR}$  and  $J^M$ , based on Monte Carlo studies, indicate that the procedure  $J^B$  has the larger expected efficiency and smaller expected size of the selected subset. The results also indicate that  $J^B$  is robust when the true distributions are not normal, that is, for some other symmetric distributions such as, the logistic, the double exponential, the tri-modal (Laplace) and the gross error model (the contaminated distributions).

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